DSC291: Machine Learning with Few Labels

Unsupervised Learning

Zhiting Hu Lecture 17, May 13, 2024



This Lecture

- Variational Inference (30mins)
- Presentation #1 (10mins):
 - Hung Nguyen, SPARF: Neural Radiance Fields from Sparse & Noisy Poses
- Presentation #2 (10mins):
 - Zhihan Chen, Efficient (Soft) Q-Learning for Text Generation with Limited Good Data

Google form for presentation questions and feedback:



• The EM algorithm:

• E-step:
$$q^{t+1} = \arg\min_{q} F\left(q, \theta^{t}\right)$$

$$= p(\mathbf{z}|\mathbf{x}, \theta^{t}) = \frac{p(\mathbf{z}, \mathbf{x}|\theta^{t})}{\sum_{z} p(\mathbf{z}, \mathbf{x}|\theta^{t})}$$

$$\circ$$
 M-step: $heta^{t+1} = rg \min_{ heta} F\left(q^{t+1}, heta^t
ight)$

$$\ell(\theta; \mathbf{x}) = \mathbb{E}_{q(\mathbf{z}|\mathbf{x})} \left[\log \frac{p(\mathbf{x}, \mathbf{z}|\theta)}{q(\mathbf{z}|\mathbf{x})} \right] + \text{KL}(q(\mathbf{z}|\mathbf{x}) || p(\mathbf{z}|\mathbf{x}, \theta))$$
$$= -F(q, \theta) + \text{KL}(q(\mathbf{z}|\mathbf{x}) || p(\mathbf{z}|\mathbf{x}, \theta))$$

- The EM algorithm:
 - \circ E-step: $q^{t+1} = \arg\min_{q} F\left(q, \theta^{t}\right)$

Intractable when model $p(\mathbf{z}, \mathbf{x}|\theta)$ is complex

$$= p(\mathbf{z}|\mathbf{x}, \theta^t) = \frac{p(\mathbf{z}, \mathbf{x}|\theta^t)}{\sum_{z} p(\mathbf{z}, \mathbf{x}|\theta^t)}$$

 \circ M-step: $heta^{t+1} = rg \min_{ heta} F\left(q^{t+1}, heta^t
ight)$

Approximate $p(\mathbf{z}|\mathbf{x}, \theta^t)$:

o find a **tractable** $q(\mathbf{z}|\mathbf{x}, \mathbf{v}^*)$ that is closest to $p(\mathbf{z}|\mathbf{x}, \theta^t)$

$$q(\mathbf{z}|\mathbf{x}, \mathbf{v}^*) = \min_{\mathbf{v}} \text{KL}(q(\mathbf{z}|\mathbf{x}, \mathbf{v}) || p(\mathbf{z}|\mathbf{x}, \theta^t))$$

$$= \min_{\mathbf{v}} F(q(\mathbf{z}|\mathbf{x},\mathbf{v}),\theta^t) + const.$$

$$\ell(\theta; \mathbf{x}) = \mathbb{E}_{q(\mathbf{z}|\mathbf{x})} \left[\log \frac{p(\mathbf{x}, \mathbf{z}|\theta)}{q(\mathbf{z}|\mathbf{x})} \right] + \text{KL}(q(\mathbf{z}|\mathbf{x}) || p(\mathbf{z}|\mathbf{x}, \theta))$$

$$= (-F(q, \theta)) + (\text{KL}(q(\mathbf{z}|\mathbf{x}) || p(\mathbf{z}|\mathbf{x}, \theta)))$$

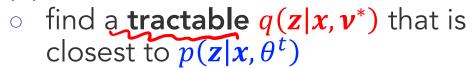


• The EM algorithm:

$$\begin{array}{c} \text{ E-step: } q^{t+1} = \underset{q}{\operatorname{arg\,min}} F\left(q,\theta^t\right) \\ \hline \text{Intractable when} \\ \operatorname{model} p(\mathbf{z},\mathbf{x}|\theta) \text{ is} \\ \operatorname{complex} \end{array} = \underbrace{p(\mathbf{z}|\mathbf{x},\theta^t)}_{q} = \underbrace{\frac{p(\mathbf{z},\mathbf{x}|\theta^t)}{\sum_{z}p(\mathbf{z},\mathbf{x}|\theta^t)}}_{q}$$

o M-step: $\theta^{t+1} = \arg\min_{q} F\left(q^{t+1}, \theta^{t}\right)$

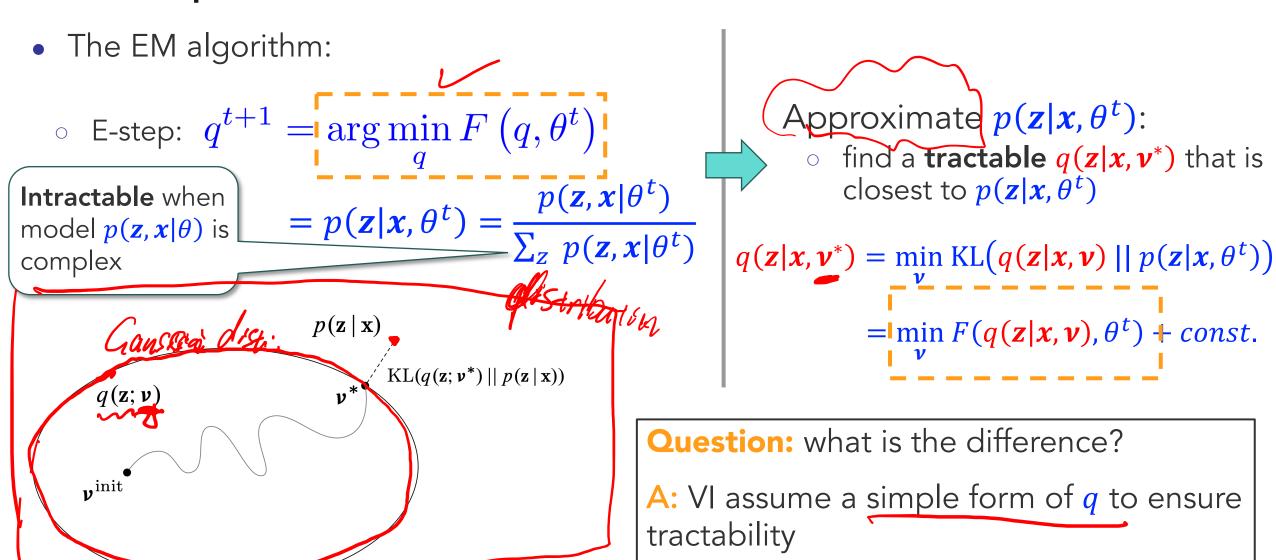
Approximate $p(\mathbf{z}|\mathbf{x}, \theta^t)$:



$$q(\mathbf{z}|\mathbf{x}, \mathbf{v}^*) = \min_{\mathbf{v}} \text{KL}(q(\mathbf{z}|\mathbf{x}, \mathbf{v}) \mid\mid p(\mathbf{z}|\mathbf{x}, \theta^t))$$
$$= \min_{\mathbf{v}} F(q(\mathbf{z}|\mathbf{x}, \mathbf{v}), \theta^t) + const.$$

Question: what is the difference?

A: VI assume a simple form of q to ensure tractability



• The EM algorithm:

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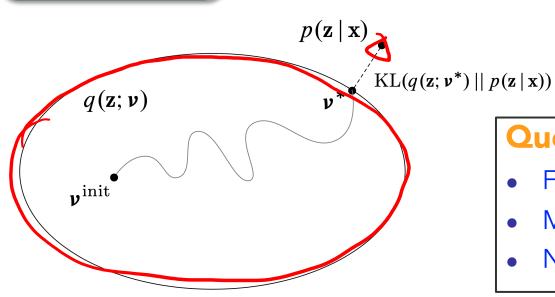
Approximate $p(\mathbf{z}|\mathbf{x}, \theta^t)$:

o find a **tractable** $q(\mathbf{z}|\mathbf{x}, \mathbf{v}^*)$ that is closest to $p(\mathbf{z}|\mathbf{x}, \theta^t)$

$$q(\mathbf{z}|\mathbf{x}, \mathbf{v}^*) = \min_{\mathbf{v}} \text{KL}(q(\mathbf{z}|\mathbf{x}, \mathbf{v}) || p(\mathbf{z}|\mathbf{x}, \theta^t))$$
$$= \min_{\mathbf{v}} F(q(\mathbf{z}|\mathbf{x}, \mathbf{v}), \theta^t) + const.$$

Question: What forms of q(z|x, y) shall we choose?

- Factorized distribution -> mean field VI
- Mixture of Gaussian distribution -> black-box VI
- Neural-based distribution -> Variational Autoencoders



Example: Mean Field Variational Inference

- A popular family of variational approximations

$$q(\mathbf{z}) = q(z_1, \dots, z_m) = \prod_{j=1}^m q(z_j)$$

- (where we omit variational parameters for ease of notation)
- We refer to $q(z_j)$, the variational approximation for a single latent variable, as a "local variational approximation"
- In the above expression, the variational approximation $q(z_j)$ over each latent variable z_i is independent

Example: Mean Field Variational Inference

- Typically, this approximation does not contain the true posterior (because the latent variables are dependent).
 - E.g.: in the (Bayesian) mixture of Gaussians model, all of the cluster assignments z_i for $i=1,\ldots,n$ are dependent on each other and on the cluster locations $\mu_{1:K}$ given data.

Example: Mean Field Variational Inference

How do we optimize the ELBO in mean field variational inference?

- Typically, we use coordinate ascent optimization.
- I.e. we optimize each latent variable's variational approximation $q(z_j)$ in turn while holding the others fixed.
 - At each iteration we get an updated "local" variational approximation.
 - And we iterate through each latent variable until convergence.

Mean Field Variational Inference with Coordinate Ascent

Recap: Bayesian mixture of Gaussians

• Treat the mean μ_k as latent variables

$$\mu_k \sim \mathcal{N}(0, \tau^2)$$
 for $k = 1, \dots, K$

• For each data i = 1, ..., n

$$z_i \sim \operatorname{Cat}(\pi)$$
.

$$x_i \sim \mathcal{N}(\mu_{z_i}, \sigma^2).$$

- We have
 - observed variables $x_{1:n}$
 - o latent variables $\mu_{1:k}$ and $z_{1:n}$
 - o parameters $\{\tau^2, \sigma^2, \pi\}$

Mean Field Variational Inference with Coordinate Ascent

Recap: Bayesian mixture of Gaussians Assume mean-field $q(\mu_{1:K}, z_{1:n}) = \prod_k q(\mu_k) \prod_i q(z_i)$ Initialize the global variational distributions $q(\mu_k)$ and parameters $\{\tau^2, \sigma^2, \pi\}$ Repeat: For each data example $i \in \{1, 2, ..., D\}$ • Update the local variational distribution $q(z_i)$ (Variational) E-step End for Update the global variational distributions $q(\mu_k)$ Update the parameters $\{\tau^2, \sigma^2, \pi\}$ → (Variational) M-step Until ELBO converges

• What if we have millions of data examples? This could be very slow.





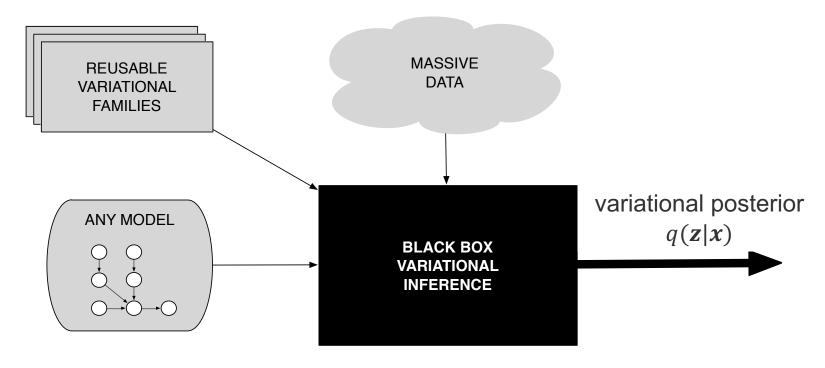
Recap: Bayesian mixture of Gaussians

Assume mean-field $q(\mu_{1:K}, z_{1:n}) = \prod_k q(\mu_k) \prod_i q(z_i)$

- Initialize the global variational distributions $q(\mu_k)$ and parameters $\{\tau^2, \sigma^2, \pi\}$
- Repeat:
 - Sample a data example $i \in \{1, 2, ..., D\}$
 - Update the local variational distribution q(z_i)
 - Update the global variational distributions $q(\mu_k)$ with natural gradient ascent
 - Update the parameters $\{\tau^2, \sigma^2, \pi\}$
- Until ELBO converges

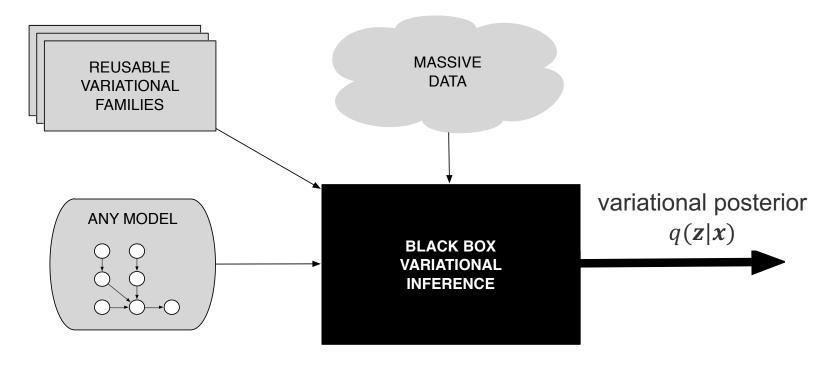
Black-box Variational Inference

- We have derived variational inference specific for Bayesian Gaussian (mixture) models
- There are innumerable models
- Can we have a solution that does not entail model-specific work?



- Easily use variational inference with any model
- Perform inference with massive data
- No mathematical work beyond specifying the model

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- Sample from q(.)
- Form noisy gradients (without model-specific computation)
- Use stochastic optimization

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- Probabilistic model: x -- observed variables, z -- latent variables
- Variational distribution $q_{\lambda}(\mathbf{z}|\mathbf{x})$ with parameters λ , e.g.,
 - Gaussian mixture distribution:
 - "A Gaussian mixture model is a universal approximator of densities, in the sense that any smooth density can be approximated with any specific nonzero amount of error by a Gaussian mixture model with enough components." (Deep Learning book, pp.65)
 - Deep neural networks
- ELBO:

$$\mathcal{L}(\lambda) = \mathbb{E}_{q(\mathbf{z}|\lambda)}[\log p(\mathbf{x}, \mathbf{z})] - \mathbb{E}_{q(\mathbf{z}|\lambda)}[\log q(\mathbf{z}|\lambda)]$$

• Want to compute the gradient w.r.t variational parameters λ

The General Problem: Computing Gradients of Expectations

• When the objective function \mathcal{L} is defined as an expectation of a (differentiable) test function $f_{\lambda}(\mathbf{z})$ w.r.t. a probability distribution $q_{\lambda}(\mathbf{z})$

$$\mathcal{L} = \mathbb{E}_{q_{\lambda}(\mathbf{z})}[f_{\lambda}(\mathbf{z})]$$

- Computing exact gradients w.r.t. the parameters λ is often unfeasible
- Need stochastic gradient estimates
 - The score function estimator (a.k.a log-derivative trick, REINFORCE)
 - The reparameterization trick (a.k.a the pathwise gradient estimator)

Computing Gradients of Expectations w/ score function

- Loss: $\mathcal{L} = \mathbb{E}_{q_{\lambda}(\mathbf{z})}[f_{\lambda}(\mathbf{z})]$
- Log-derivative trick: $\nabla_{\lambda} q_{\lambda} = q_{\lambda} \nabla_{\lambda} \log q_{\lambda}$
- Gradient w.r.t. λ:

$$\nabla_{\lambda} \mathcal{L} = \mathbb{E}_{q_{\lambda}(\mathbf{z})} [f_{\lambda}(\mathbf{z}) \nabla_{\lambda} \log q_{\lambda}(\mathbf{z}) + \nabla_{\lambda} f_{\lambda}(\mathbf{z})]$$

- o score function: the gradient of the log of a probability distribution
- ullet Compute noisy unbiased gradients with Monte Carlo samples from q_λ

$$\nabla_{\lambda} \mathcal{L} \approx \frac{1}{S} \sum_{s=1}^{S} f_{\lambda}(\mathbf{z}_{s}) \nabla_{\lambda} \log q_{\lambda}(\mathbf{z}_{s}) + \nabla_{\lambda} f_{\lambda}(\mathbf{z}_{s})$$
 where $\mathbf{z}_{s} \sim q_{\lambda}(\mathbf{z})$

- Pros: generally applicable to any distribution $q(z|\lambda)$
- Cons: empirically has high variance → slow convergence
 - To reduce variance: Rao-Blackwellization, control variates, importance sampling, ...

Computing Gradients of Expectations w/ reparametrization trick

- Loss: $\mathcal{L} = \mathbb{E}_{q_{\lambda}(\mathbf{z})}[f_{\lambda}(\mathbf{z})]$
- Assume that we can express the distribution $q_{\lambda}(z)$ with a transformation

$$\begin{array}{l}
\epsilon \sim s(\epsilon) \\
z = t(\epsilon, \lambda)
\end{array} \iff z \sim q(z|\lambda)$$

E.g.,

$$\begin{array}{l} \epsilon \sim Normal(0,1) \\ z = \epsilon \sigma + \mu \end{array} \iff z \sim Normal(\mu, \sigma^2)$$

Reparameterization gradient

$$\mathcal{L} = \mathbb{E}_{\epsilon \sim S(\epsilon)}[f_{\lambda}(\mathbf{z}(\epsilon, \lambda))]$$

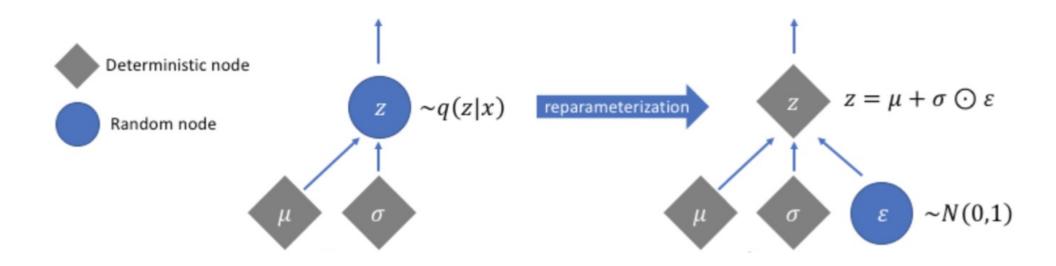
$$\nabla_{\lambda} \mathcal{L} = \mathbb{E}_{\epsilon \sim S(\epsilon)}[\nabla_{\mathbf{z}} f_{\lambda}(\mathbf{z}) \nabla_{\lambda} t(\epsilon, \lambda)]$$

- Pros: empirically, lower variance of the gradient estimate
- Cons: Not all distributions can be reparameterized

Reparameterization trick

• Reparametrizing Gaussian distribution

$$\begin{array}{l} \epsilon \sim Normal(0,1) \\ z = \epsilon \sigma + \mu \end{array} \iff z \sim Normal(\mu, \sigma^2)$$



[Courtesy: Tansey, 2016]

Reparameterization trick

• Reparametrizing Gaussian distribution

$$\begin{array}{l} \epsilon \sim Normal(0,1) \\ z = \epsilon \sigma + \mu \end{array} \iff z \sim Normal(\mu, \sigma^2)$$

- Other reparameterizable distributions: $\epsilon \sim Uniform(\epsilon)$ • Tractable inverse CDF F^{-1} : $z = F^{-1}(\epsilon)$ $\Leftrightarrow z \sim q(z)$
 - Exponential, Cauchy, Logistic, Rayleigh, Pareto, Weibull, Reciprocal, Gompertz, Gumbel, Erlang
 - Location-scale:
 - Laplace, Elliptical, Student's t, Logistic, Uniform, Triangular, Gaussian
 - Composition:
 - Log-Normal (exponentiated normal) Gamma (sum of exponentials) Dirichlet (sum of Gammas) Beta, Chi-Squared, F

[Courtesy: Tansey, 2016]

Computing Gradients of Expectations: Summary

- Loss: $\mathcal{L} = \mathbb{E}_{q_{\lambda}(\mathbf{z})}[f_{\lambda}(\mathbf{z})]$
- Score gradient

$$\nabla_{\lambda} \mathcal{L} = \mathbb{E}_{q_{\lambda}(\mathbf{z})} [f_{\lambda}(\mathbf{z}) \nabla_{\lambda} \log q_{\lambda}(\mathbf{z}) + \nabla_{\lambda} f_{\lambda}(\mathbf{z})]$$

- Pros: generally applicable to any distribution $q(z|\lambda)$
- Cons: empirically has high variance → slow convergence
- Reparameterization gradient

$$\nabla_{\lambda} \mathcal{L} = \mathbb{E}_{\epsilon \sim s(\epsilon)} [\nabla_{\mathbf{z}} f_{\lambda}(\mathbf{z}) \nabla_{\lambda} t(\epsilon, \lambda)]$$

- o Pros: empirically, lower variance of the gradient estimate
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- Variational distribution $q_{\lambda}(\mathbf{z}|\mathbf{x})$ with parameters λ , e.g.,
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 - Deep neural networks $\mathcal{L}(\lambda) \triangleq \mathrm{E}_{q_{\lambda}(z)}[\log p(x,z) \log q(z)].$
- ELBO:

$$\mathcal{L}(\lambda) = \mathbb{E}_{q(\mathbf{z}|\lambda)}[\log p(\mathbf{x}, \mathbf{z})] - \mathbb{E}_{q(\mathbf{z}|\lambda)}[\log q(\mathbf{z}|\lambda)]$$

• Want to compute the gradient w.r.t variational parameters λ

BBVI with the score gradient

• ELBO:

$$\mathcal{L}(\lambda) = \mathbb{E}_{q(\mathbf{z}|\lambda)}[\log p(\mathbf{x}, \mathbf{z})] - \mathbb{E}_{q(\mathbf{z}|\lambda)}[\log q(\mathbf{z}|\lambda)]$$

• Gradient w.r.t. λ (using the log-derivative trick)

$$\nabla_{\lambda} \mathcal{L} = \mathrm{E}_{q}[\nabla_{\lambda} \log q(z|\lambda)(\log p(x,z) - \log q(z|\lambda))]$$

 Compute noisy unbiased gradients of the ELBO with Monte Carlo samples from the variational distribution

$$abla_{\lambda} \mathcal{L} pprox rac{1}{S} \sum_{s=1}^{S}
abla_{\lambda} \log q(z_s | \lambda) (\log p(x, z_s) - \log q(z_s | \lambda)),$$
where $z_s \sim q(z | \lambda)$.

BBVI with the reparameterization gradient

• ELBO:

$$\mathcal{L}(\lambda) = \mathbb{E}_{q(\mathbf{z}|\lambda)}[\log p(\mathbf{x}, \mathbf{z})] - \mathbb{E}_{q(\mathbf{z}|\lambda)}[\log q(\mathbf{z}|\lambda)]$$

Gradient w.r.t. λ

$$\begin{array}{l} \epsilon \sim s(\epsilon) \\ z = t(\epsilon, \lambda) \end{array} \iff z \sim q(z|\lambda)$$

$$\nabla_{\lambda} \mathcal{L} = \mathbb{E}_{\epsilon \sim s(\epsilon)} \left[\nabla_z [\log p(x, z) - \log q(z)] \nabla_{\lambda} t(\epsilon, \lambda) \right]$$

Questions?