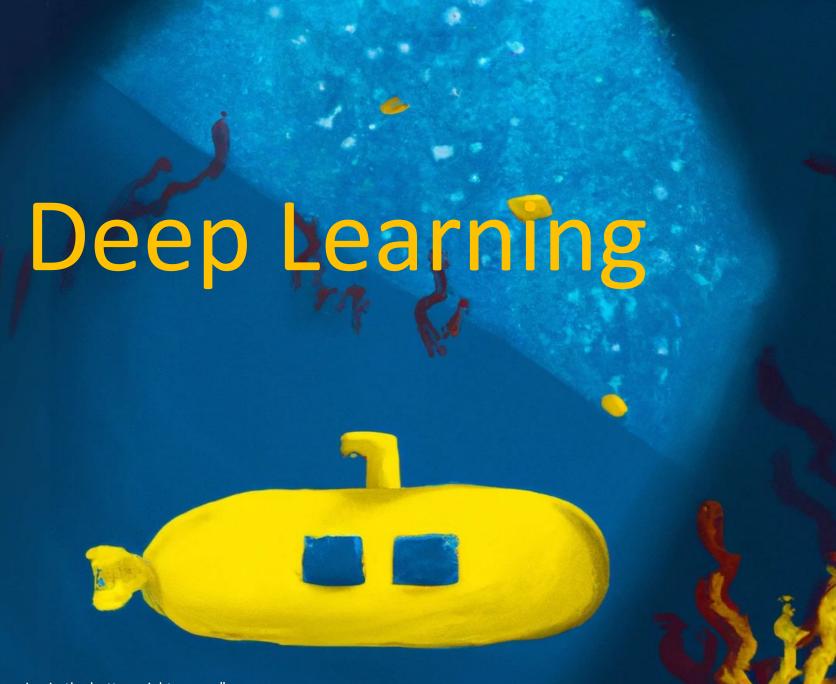
CSCI 1470/2470 Spring 2023

Ritambhara Singh

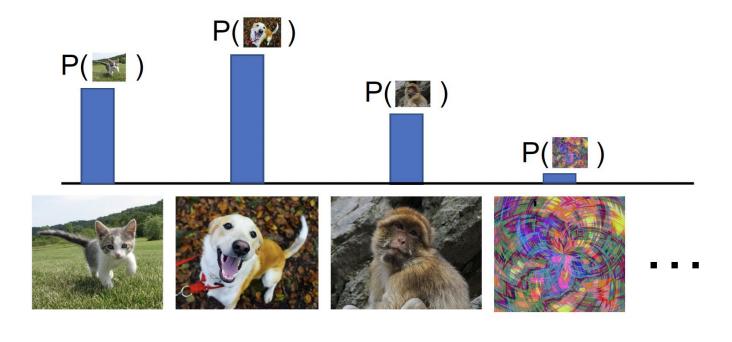
April 07, 2023 Friday



### Review: Discriminative v/s Generative models

Discriminative Model: Learn a probability distribution p(y|x)

**Generative Model**: Learn a probability distribution p(x)



- Generative model: All possible images compete with each other for probability mass
- Model can "reject" unreasonable inputs by assigning them small values

Credit: UMich EECS498

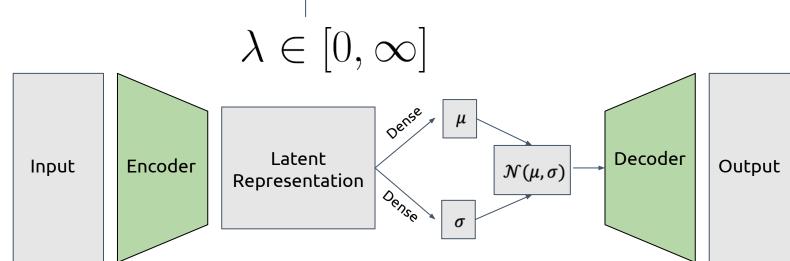
### Review: Weighted Combination of Losses

 $L_1$  = loss associated with producing output similar to input

 ${\cal L}_2$  = loss associated with producing output with some variation to input

$$L = L_1 + \lambda L_2$$

**Total Loss:** 



# Today's goal – continue to learn about variational autoencoders (VAEs)

- (1) VAE Loss KL Divergence
- (2) Reparameterization trick
- (3) Conditional VAE

### VAE Losses, Defined

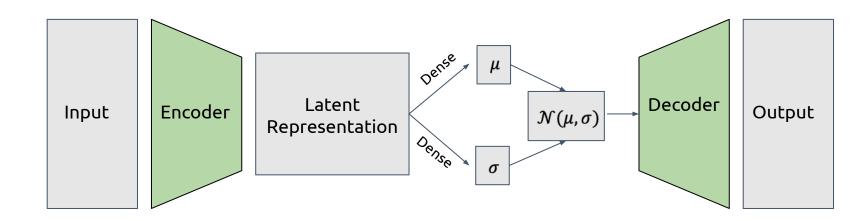
We have seen  ${\cal L}_1$  before: this is just the autoencoder reconstruction loss

$$L_1(x,\hat{x}) = ||x - \hat{x}||_2^2$$

But with  $L_2$ , it's not so clear. How do we measure how much variation our output would have?

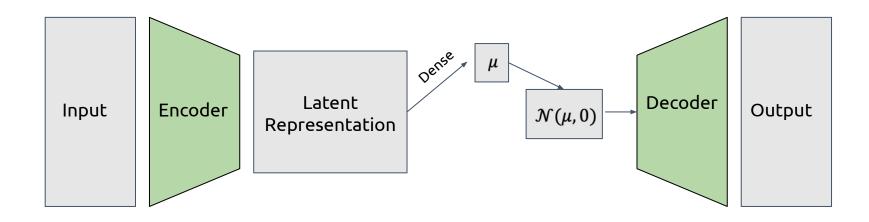
$$L_2(??,??) = ??$$

- ullet To get variation, we definitely need a loss that encourages  $\sigma>0$ 
  - If we don't do this,  $L_1$  will drive  $\sigma$  to zero in an effort to produce the best reconstructions

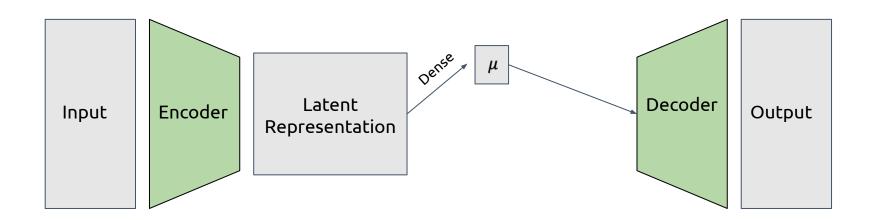


- To get variation, we definitely need a loss that encourages  $\sigma>0$ 
  - If we don't do this,  $L_1$  will drive  $\sigma$  to zero in an effort to produce the best reconstructions

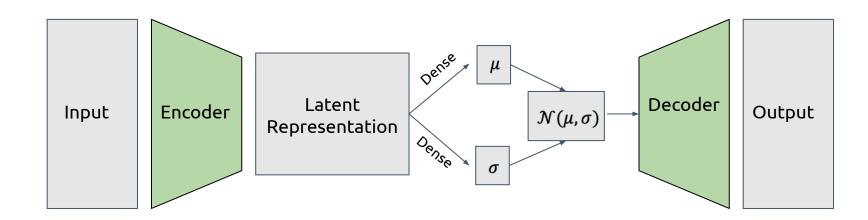
What's the issue here?



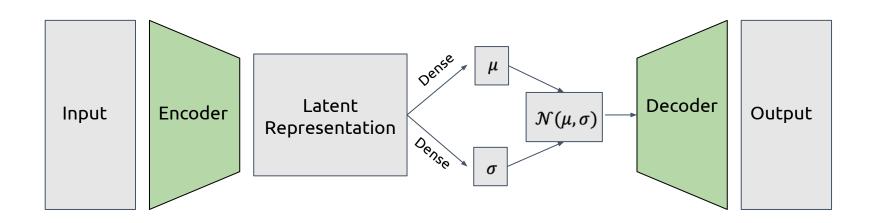
- To get variation, we definitely need a loss that encourages  $\sigma>0$ 
  - If we don't do this,  $L_1$  will drive  $\sigma$  to zero in an effort to produce the best reconstructions
  - Behaves the same as a regular autoencoder!



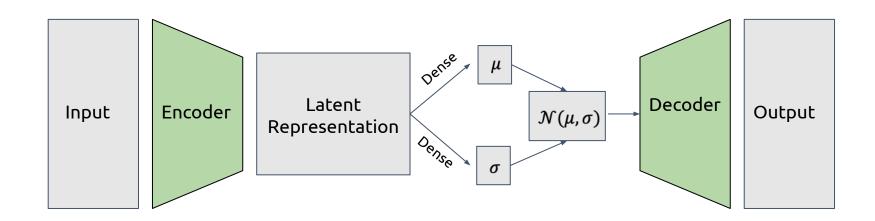
- To get variation, we definitely need a loss that encourages  $\sigma>0$ 
  - If we don't do this,  $L_1$  will drive  $\sigma$  to zero in an effort to produce the best reconstructions
- But how big should we encourage  $\sigma$  to be?
- And for that matter, what we do about  $\mu$ ?



- The idea: make  $\mathcal{N}(\mu, \sigma)$  close to  $\mathcal{N}(0, 1)$ 
  - Obviously, we can't perfectly satisfy this for every input (otherwise every input would produce the same set of outputs → terrible reconstruction!)
  - But, we'll see later that having some light pressure to make  $\mathcal{N}(\mu, \sigma)$  close to  $\mathcal{N}(0, 1)$  will have some beneficial properties



- Wait...but **how** do we make  $\mathcal{N}(\mu, \sigma)$  close to  $\mathcal{N}(0, 1)$ ?
- More generally: how do measure the difference between two probability distributions?



Measures the difference between any two probability distributions

$$D_{KL}(P||Q) = \int_{-\infty}^{\infty} \frac{p(x)}{p(x)} \log\left(\frac{p(x)}{q(x)}\right) dx$$

### What this says:

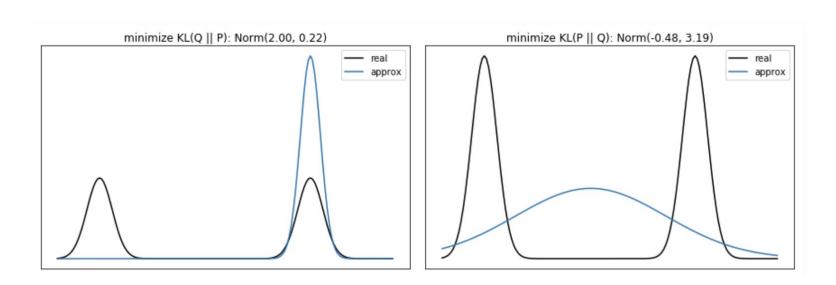
- "Everywhere that p has probability density..."
- "...the difference between p and q should be small"
  - Difference in log probabilities (remember that  $\log\left(\frac{a}{b}\right) = \log(a) \log(b)$ )

More on KL Divergence:

Measures the difference between any two probability distributions

$$D_{KL}(P||Q) = \int_{-\infty}^{\infty} p(x) \log \left(\frac{p(x)}{q(x)}\right) dx$$

• Note that this is not symmetric:  $D_{KL}(P||Q) \neq D_{KL}(Q||P)$ 



- Expensive to compute, in general (no closed form, have to numerically approximate the integral)
- But! There is a closed form for Gaussians:

$$D_{KL}(\mathcal{N}(\mu, \sigma^2) || \mathcal{N}(0, 1)) = \frac{1}{2} \sum_{i=1}^{k} (\mu_i^2 + \sigma_i^2 - \ln \sigma_i^2 - 1)$$

k is the dimensionality of  $m{\mu}$  and  $m{\sigma}$  (e.g. k = 100 when  $\mu \in \mathbb{R}^{100}$  )

We won't derive the equation above, but let's convince ourselves it behaves how we expect it to behave

- Expensive to compute, in general (no closed form, have to numerically approximate the integral)
- But! There is a closed form for Gaussians:

$$D_{KL}(\mathcal{N}(\mu, \sigma^2) || \mathcal{N}(0, 1)) = \frac{1}{2} \sum_{i=1}^{k} (\mu_i^2 + \sigma_i^2 - \ln \sigma_i^2 - 1)$$

Derive the expression for (1)  $\sigma$ =1 and (2)  $\mu$ =0

### KL Divergence for Two Gaussians

$$D_{KL}(\mathcal{N}(\mu, \sigma^2) || \mathcal{N}(0, 1)) = \frac{1}{2} \sum_{i=1}^{k} (\mu_i^2 + \sigma_i^2 - \ln \sigma_i^2 - 1)$$

Let's take the case  $\sigma = 1$ 

$$D_{KL}(\mathcal{N}(\mu, 1)||\mathcal{N}(0, 1)) = \frac{1}{2} \sum_{i=1}^{k} (\mu_i^2 + 1^2 - \ln(1) - 1)$$
$$= \frac{1}{2} \sum_{i=1}^{k} \mu_i^2$$

The expression is minimized  $\mu = 0$  (which is what we want!)

### KL Divergence for Two Gaussians

$$D_{KL}(\mathcal{N}(\mu, \sigma^2) || \mathcal{N}(0, 1)) = \frac{1}{2} \sum_{i=1}^{k} (\mu_i^2 + \sigma_i^2 - \ln \sigma_i^2 - 1)$$

Let's take the case  $\mu = 0$ 

$$D_{KL}(\mathcal{N}(0,\sigma^2)||\mathcal{N}(0,1)) = \frac{1}{2} \sum_{i=1}^{k} (\sigma_i^2 - \ln(\sigma_i^2) - 1)$$

This expression is minimized when  $\sigma = 1$  (which is also what we want!)

### The Final VAE Loss Function

We now have all the tools necessary to construct our loss function.

$$L = L_1 + \lambda L_2 \qquad \qquad \lambda \in [0, \infty]$$

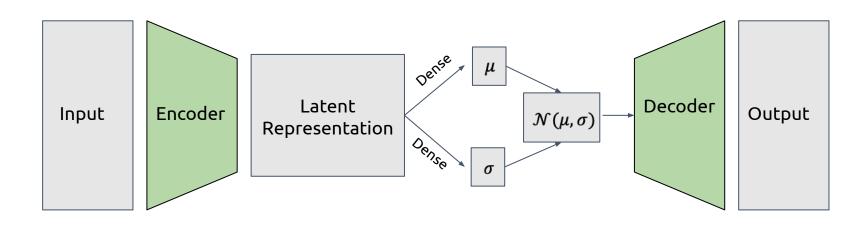
Which turns into this:

$$L = ||x - \hat{x}||_2^2 + \lambda D_{KL}(\mathcal{N}(\mu, \sigma), \mathcal{N}(0, 1))|$$



### Putting it all together

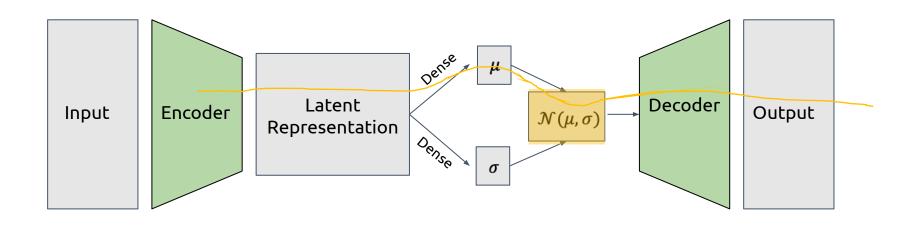
$$L = ||x - \hat{x}||_2^2 + \lambda D_{KL}(\mathcal{N}(\mu, \sigma), \mathcal{N}(0, 1))|$$



### Ah, but there's a catch:

Can anyone guess?

- To update the weights of the encoder, we have to backprop through a random sampling operation
- Sampling a random value seems not differentiable...



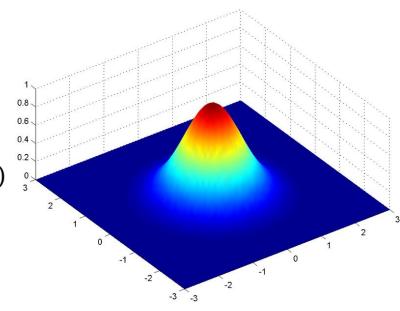
# Remember our sampling strategy for Gaussian?

•

The Gaussian Distribution

• 
$$p(x \mid \mu, \sigma) = \mathcal{N}(\mu, \sigma)(x) = \frac{1}{\sqrt{2\pi\sigma^2}} e^{-\frac{(x-\mu)^2}{2\sigma^2}}$$

- Sampling:
  - Sample from the unit normal distribution  $\rightarrow r \sim \mathcal{N}(0,1)$
  - Return  $\mu + r\sigma$



### The Reparameterization Trick

A nice property of Gaussian distributions: if we sample  $z \sim \mathcal{N}(\mu, \sigma)$  we can rewrite it as:

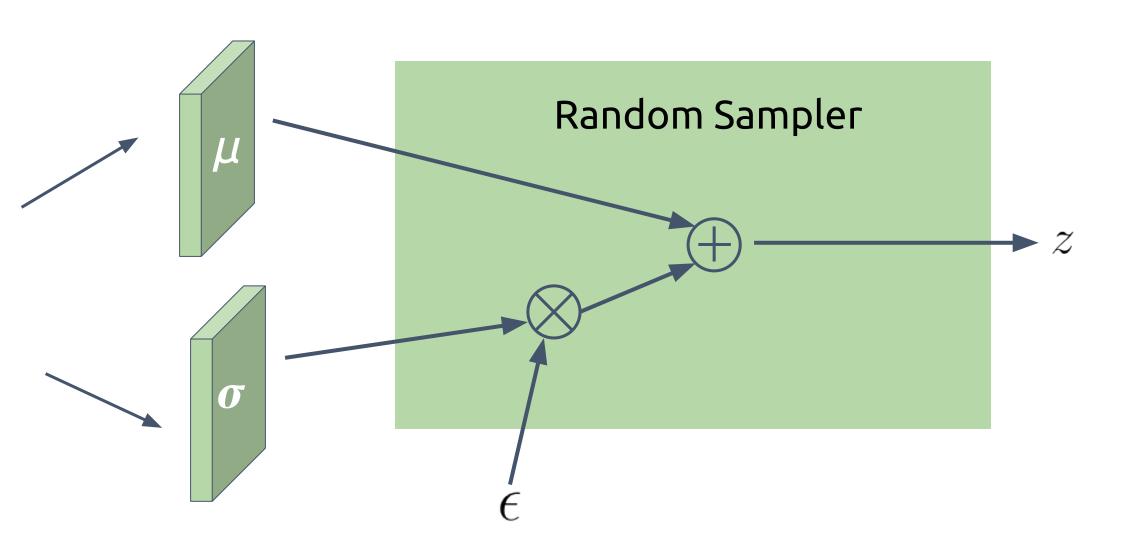
$$z = \mu + \epsilon \cdot \sigma$$

Where

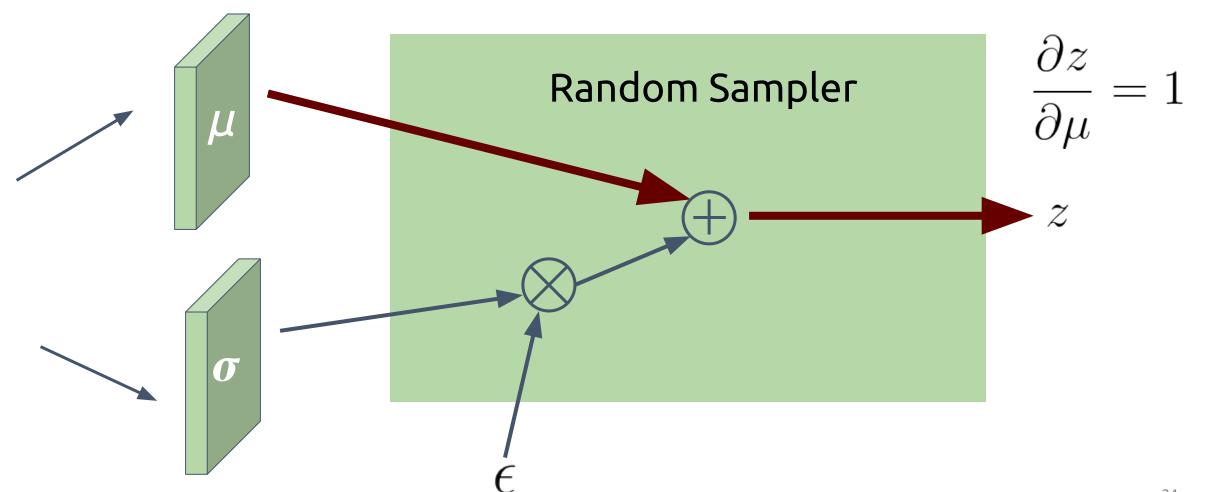
$$\epsilon \sim \mathcal{N}(0,1)$$

- The random sampling no longer depends on learnable parameters
- This allows us to do backpropagation

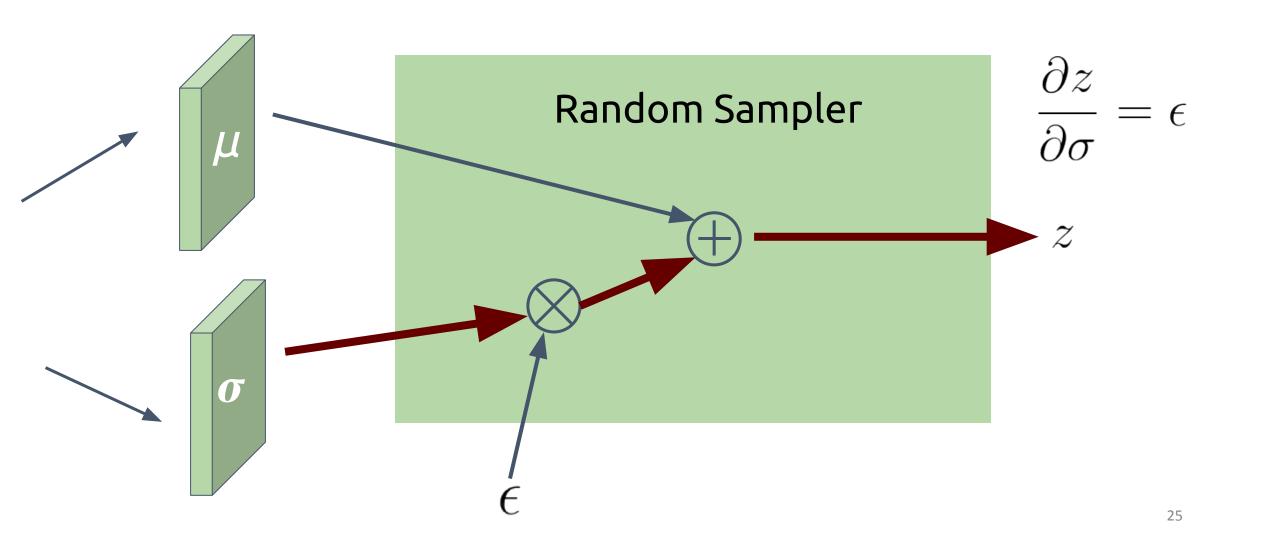
### Random Sampler with Reparameterization Trick



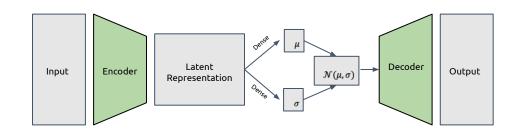
### Random Sampler with Reparameterization Trick



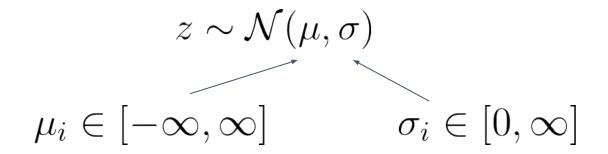
### Random Sampler with Reparameterization Trick



### One more practical detail

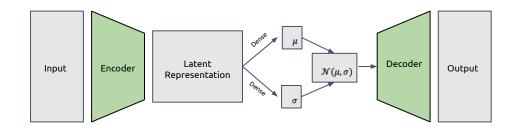


Let's again consider our sampling operation

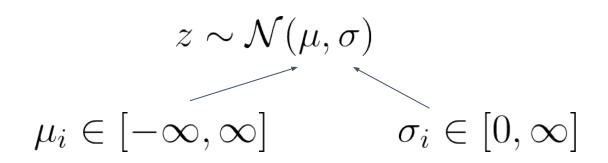


- Nothing prevents the neural network from outputting negative values for the standard deviation.
- Instead of predicting  $\sigma$ , we will instead predict  $\log(\sigma^2)$  . This ensures that every  $\sigma_i \in [0,\infty]$ 
  - ullet i.e. just treat the output of the Dense layer as if it is  $\log(\sigma^2)$

### One more practical detail



Let's again consider our sampling operation



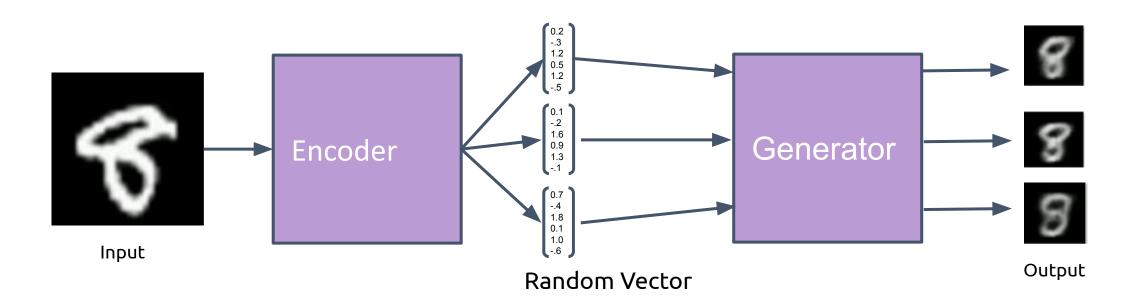


- Instead of predicting  $\sigma$ , we will instead predict  $\log(\sigma^2)$  . This ensures that every  $\sigma_i \in [0,\infty]$ 
  - ullet i.e. just treat the output of the Dense layer as if it is  $\log(\sigma^2)$

$$D_{KL}(\mathcal{N}(\mu, \sigma^2) || \mathcal{N}(0, 1)) = \frac{1}{2} \sum_{i=1}^{k} (\mu_i^2 + \sigma_i^2 - \ln \sigma_i^2 - 1)$$

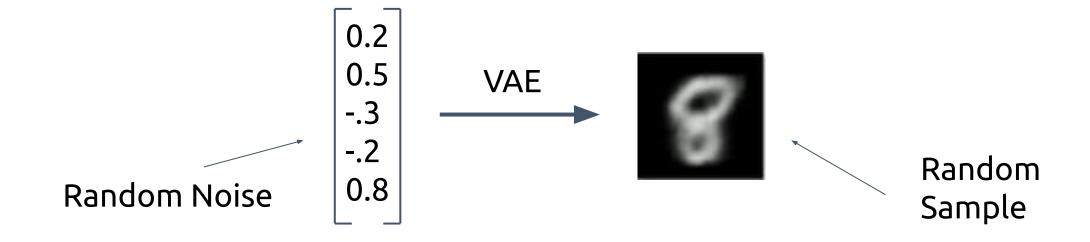
### Sampling from a VAE

• We can use a trained VAE to generate random variants of an input data point...



### Sampling from a VAE

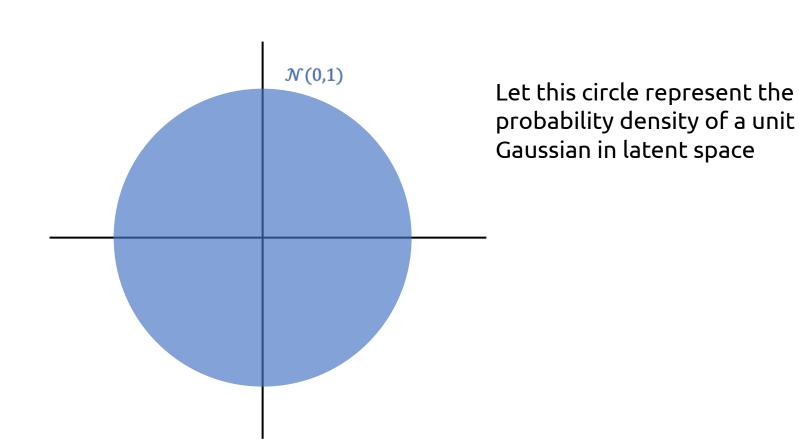
... But ultimately, we want to draw random samples from a VAE



#### How can we do this?

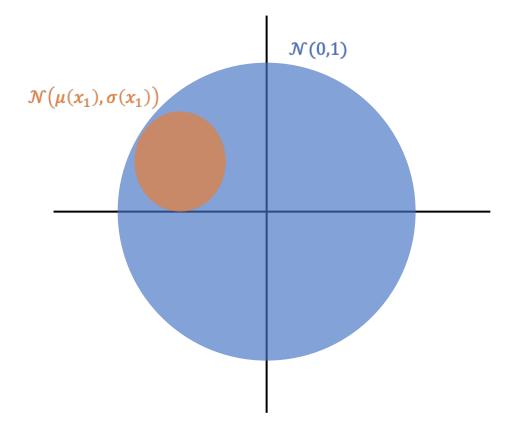
This is where our particular choice of training loss will pay off

## Encoding different points into latent space



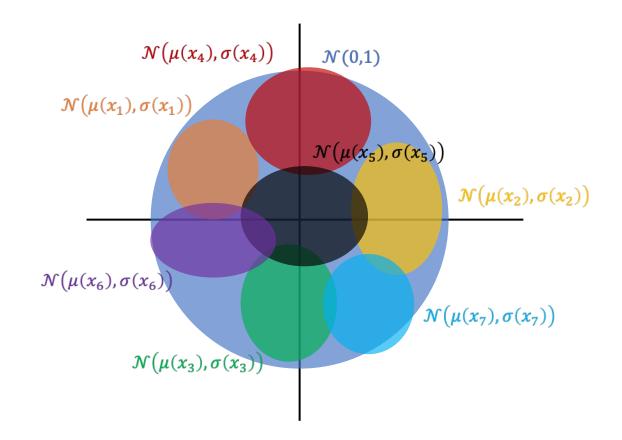
### Encoding different points into latent space

Let this circle represent the probability density of the  $\mathcal{N}(\mu,\sigma)$  distribution that the encoder predicts given an input data point  $x_1$ 



### Encoding different points into latent space

$$L = ||x - \hat{x}||_{2}^{2} + \lambda D_{KL}(\mathcal{N}(\mu, \sigma), \mathcal{N}(0, 1))|$$

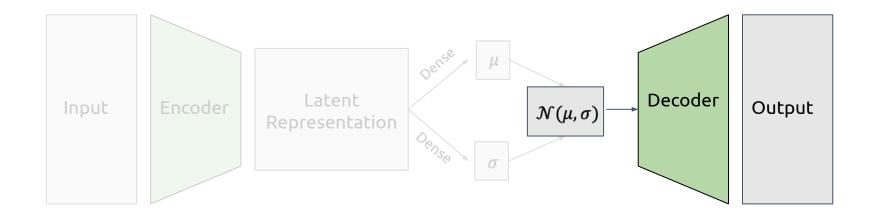


Because of our KL divergence loss, the  $\mathcal{N}(\mu, \sigma)$  for any input data point has to be somewhat similar to  $\mathcal{N}(0,1)$ 

So, if we sample a point from  $\mathcal{N}(0,1)$ , it is very likely to fall within one of these encoded

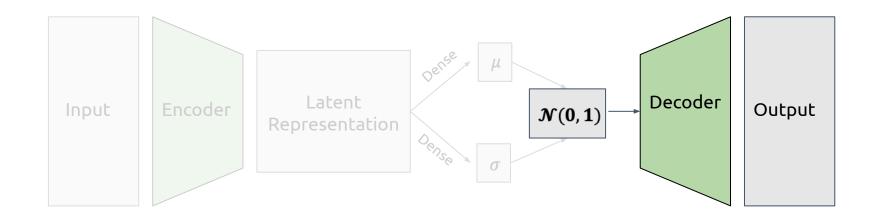
#### So what do we do?

## Sampling from a VAE



• Discard this part of the network...

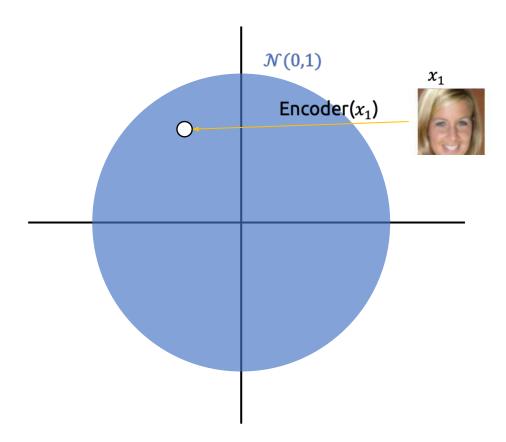
### Sampling from a VAE



- Discard this part of the network...
- ...and set  $(\mu, \sigma) = (0, 1)$

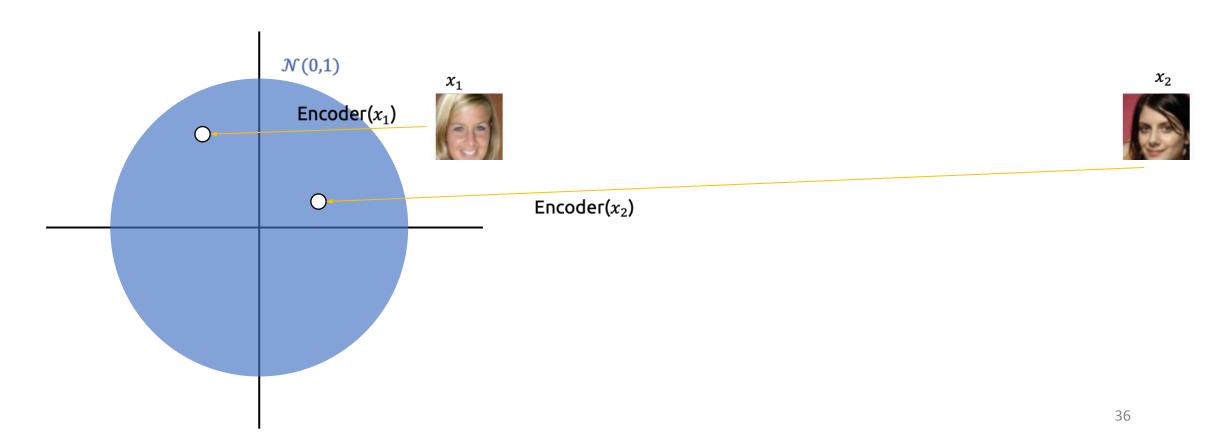
### Latent Space Interpolation

• Trace a linear path between two points in latent space, put all points along the path into the decoder



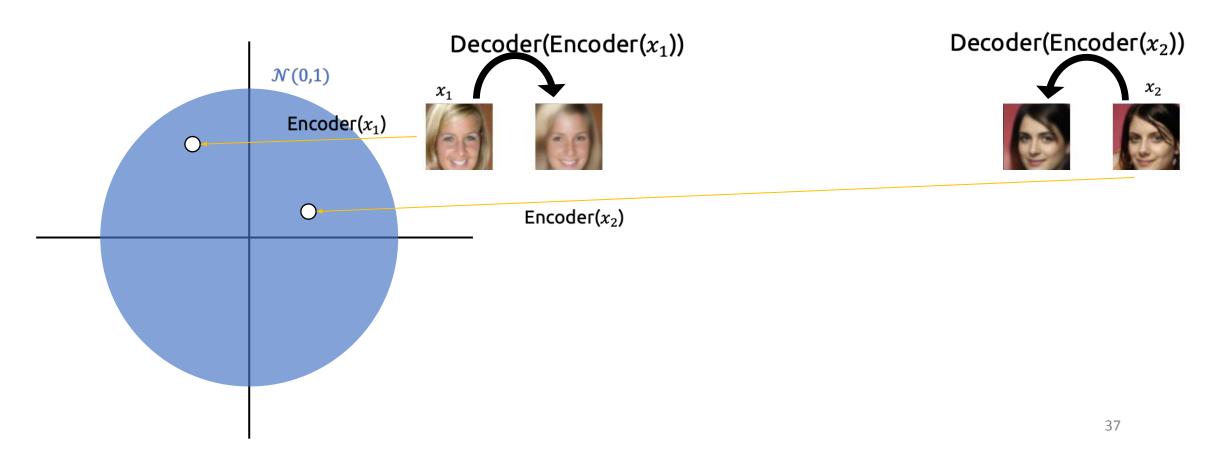
### Latent Space Interpolation

• Trace a linear path between two points in latent space, put all points along the path into the decoder



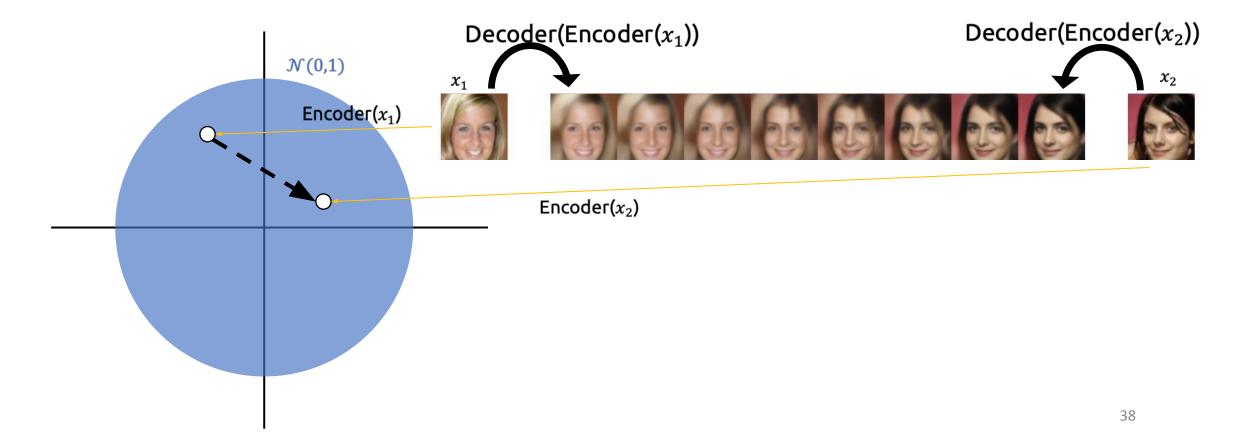
### Latent Space Interpolation

• Trace a linear path between two points in latent space, put all points along the path into the decoder



### Latent Space Interpolation

• Trace a linear path between two points in latent space, put all points along the path into the decoder





### Latent Space Interpolation

- Can also try it with a regular autoencoder
  - Doesn't work as well
  - Why not?
  - The KL divergence loss regularizes the shape of the latent space. Without it, a regular autoencoder might have "empty" pockets of latent space

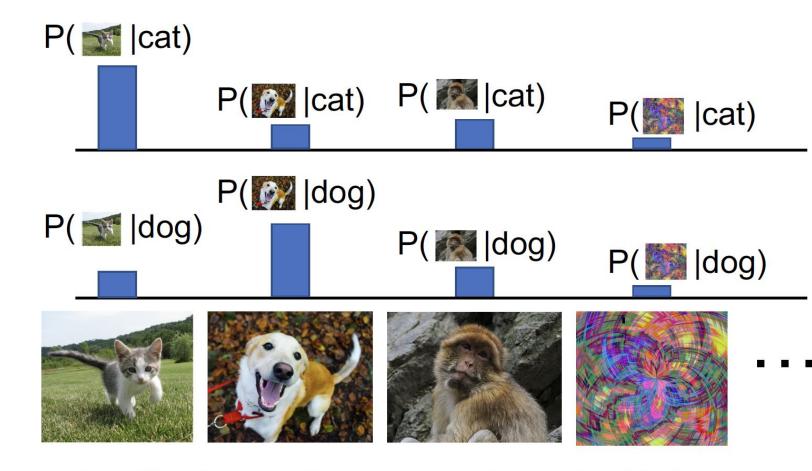


### Discriminative vs Generative Models

Discriminative Model: Learn a probability distribution p(y|x)

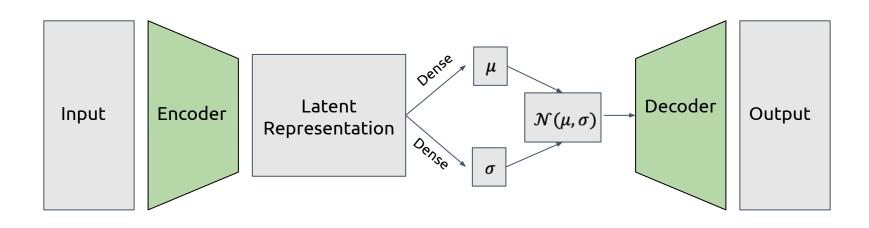
Generative Model: Learn a probability distribution p(x)

**Conditional Generative Model:** Learn p(x|y)

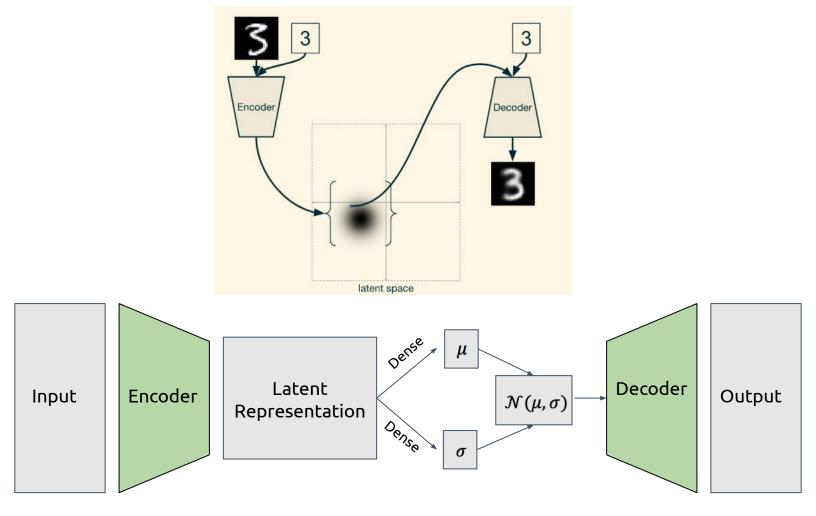


Conditional Generative Model: Each possible label induces a competition among all images

### **Conditional VAE**



### **Conditional VAE**



# VAE output

#### Input



VAE reconstruction



https://towardsdatascience.com/what-the-heck-are-vae-gans-17b86023588a

What's the issue here?

Why?

### Why are VAE samples blurry?

- Our reconstruction loss is the culprit
- Mean Square Error (MSE) loss looks at each pixel in isolation
- If no pixel is too far from its target value, the loss won't be too bad
- Individual pixels look OK, but larger-scale features in the image aren't recognizable

#### • Solutions?

Let's choose a different reconstruction loss!

#### Input



VAE reconstruction



https://towardsdatascience.com/what-the-heck-are-vae-gans-17b86023588a

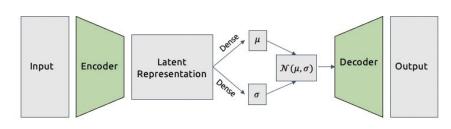
# Recap

Variational Autoencoders (VAEs)



Reparameterization Trick

**Conditional VAEs** 









VAE reconstruction



https://towardsdatascience.com/wha t-the-heck-are-vae-gans-17b860235 88a

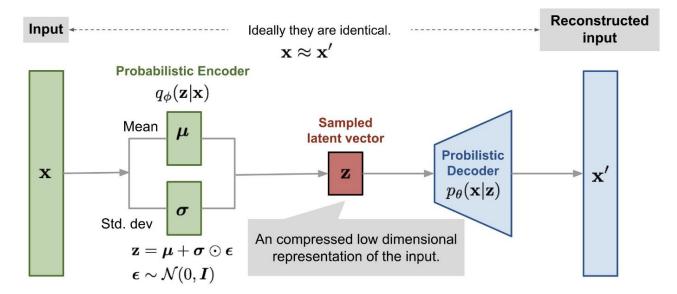
Extra Material: Deriving the VAE loss

Full derivation here - <a href="https://arxiv.org/pdf/1907.08956.pdf">https://arxiv.org/pdf/1907.08956.pdf</a>

Unfortunately, z is unknown, so we need to **marginalize** over all possible z:

$$p_{\theta}(x) = \int p_{\theta}(x, z) dz = \int p_{\theta}(x|z) p_{\theta}(z) dz$$

### Problem: Impossible to integrate over all z!



How to train this model?

Basic idea: maximize likelihood of data

compute with decoder network we assume Gaussian prior

<sup>\*</sup>Marginalization is a method that requires summing over the possible values of one variable to determine the marginal contribution of another

compute with decoder

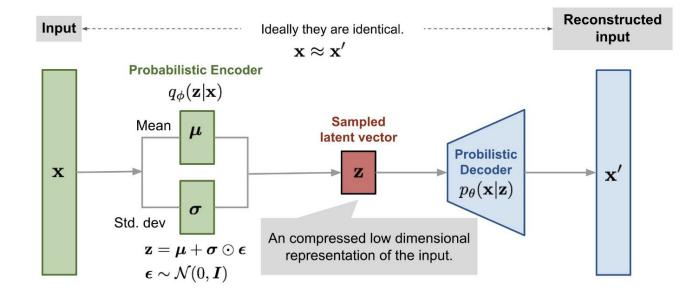
we assume Gaussian prior

$$p_{\theta}(x) = \frac{p_{\theta}(x \mid z)p_{\theta}(z)}{p_{\theta}(z \mid x)}$$

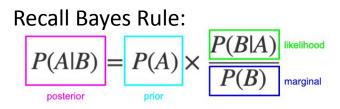
$$\approx \frac{p_{\theta}(x \mid z)p_{\theta}(z)}{q_{\phi}(z \mid x)}$$

Train an encoder that learns

$$q_{\phi}(z \mid x) \approx p_{\theta}(z \mid x)$$



**Idea**: Jointly train both encoder and decoder to maximize  $p_{\theta}(x)$ !



$$p_{\theta}(x) = \frac{p_{\theta}(x \mid z)p_{\theta}(z)}{p_{\theta}(z \mid x)}$$

Bayes' Rule

$$\log p_{\theta}(x) = \log \frac{p_{\theta}(x \mid z)p(z)}{p_{\theta}(z \mid x)}$$

Take log on each sides

$$= \log \frac{p_{\theta}(x|z)p(z)q_{\phi}(z|x)}{p_{\theta}(z|x)q_{\phi}(z|x)}$$

Multiply top and bottom by  $q_{\Phi}(z|x)$ 

$$= \log p_{\theta}(x|z) - \log \frac{q_{\phi}(z|x)}{p(z)} + \log \frac{q_{\phi}(z|x)}{p_{\theta}(z|x)}$$

Split up using rules for logarithms

$$\log p_{\theta}(x) = \log p_{\theta}(x|z) - \log \frac{q_{\phi}(z|x)}{p(z)} + \log \frac{q_{\phi}(z|x)}{p_{\theta}(z|x)}$$

We want to maximize the likelihood of the distribution p(x)

So, we reframe the likelihood function by wrapping in expectation w.r.t. z

$$= E_{z}[\log p_{\theta}(x|z)] - E_{z}\left[\log \frac{q_{\phi}(z|x)}{p(z)}\right] + E_{z}\left[\log \frac{q_{\phi}(z|x)}{p_{\theta}(z|x)}\right]$$

$$\downarrow \qquad \qquad \downarrow$$

$$= E_{z \sim q_{\phi}(z|x)}[\log p_{\theta}(x|z)] - D_{KL}\left(q_{\phi}(z|x), p(z)\right) + D_{KL}(q_{\phi}(z|x), p_{\theta}(z|x))$$

 $\log p_{\theta}(x) = E_{z \sim q_{\phi}(z|x)}[\log p_{\theta}(x)]$ 

doesn't depend on z

$$D_{ ext{KL}}(P \parallel Q) = \sum_{x \in \mathcal{X}} P(x) \log igg(rac{P(x)}{Q(x)}igg)$$

Data reconstruction by the decoder

KL divergence between prior, and samples from the encoder network

KL is >= 0, so dropping this term gives a **lower bound!** 

$$\log p_{\theta}(x) \geq E_{z \sim q_{\phi}(z|x)}[\log p_{\theta}(x|z)] - D_{KL}\left(q_{\phi}(z|x), p(z)\right) \quad \text{Variational Lower Bound}$$

Maximum Likelihood Estimation:

$$\log p_{\theta}(x) \ge E_{z \sim q_{\phi}(z|x)}[\log p_{\theta}(x|z)] - D_{KL}\left(q_{\phi}(z|x), p(z)\right)$$

Loss:

$$-E_{z\sim q_{\phi}(z|x)}[\log p_{\theta}(x|z)] + D_{KL}\left(q_{\phi}(z|x), p(z)\right)$$

$$L = ||x - \hat{x}||_2^2 + \lambda D_{KL}(\mathcal{N}(\mu, \sigma), \mathcal{N}(0, 1))|$$

See Deep Learning Book (Section 5.5)