# Halvtid presentation

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HIS

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#### Autoencoder

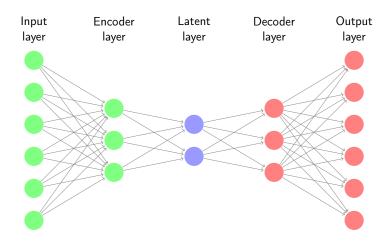
#### **Basics**

- "been part of the historical landscape of of neural networks for decades" [3]
- ▶ Hidden layers h reassemble a function f(X) which transform their input X to a encoded representation z so that another reconstruction function g(z) is able to produce  $g(z) \approx X$ .

## Usage

- Denoising
- Dimensionality reduction
- ► Feature learning
- ► Generative learning

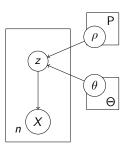
## Autoencoder



#### Method

Traditionally unsupervised variational Bayesian method

### Sampling process



- Model parameter space
- $\rho$  Probability distribution P(z)
- z Latent variable sampled n times

### Bayes theorem

$$P(A \mid B) = \frac{P(B \mid A)P(A)}{P(B)} \tag{1}$$

### Objective

- ▶ With every example X from a set  $\mathcal{X}$  we aim to generate a space z to recreate X.
- ▶ z gets sampled from  $\mathcal{Z}$  according to  $\rho \sim P(z)$ .
- ▶  $f(z;\theta), f: \mathcal{Z} \times \Theta \to \mathcal{X}$  helps to optimize  $\theta$
- maximize:

$$P(x) = \int P(X \mid z; \theta) P(z) dz$$
 (2)

▶ Where  $f(z; \theta)$  gets replaced by  $P(X \mid z; \theta)$  to enable maximum likelihood framework

### Objective

Based on some input sample values of z likely to produce X.

#### Method

Therefor a function  $Q(z \mid X)$  is introduced to generate z's likely to produce X.

This space of z's should be smaller than that one under the prior P(X)

#### Optimisation

The framework introduced by Kingma and Welling [4] does the magic. It allows us to construct a differentiable estimator:

$$\int Q(z \mid X) f(z) dz \tag{3}$$

### Kullback Leibler Divergence

Allows to compare probability distributions

P, Q probability distribution functions

$$\mathcal{D}(P \parallel Q) = \mathcal{KL} = \sum_{x \in X} P(x) \cdot \log \frac{P(x)}{Q(x)}$$
 (4)

# Helpers

### Marginal likelihood of individual data points

(Likelihood function integrated over parameter space)

$$\log P\left(X^{i}\right) = \mathcal{D}\left(Q_{\phi}\left(z\mid X^{i}\right) \parallel P_{\theta}\left(z\mid X^{i}\right)\right) + \mathcal{L}\left(\theta, \phi, X^{i}\right) \quad (5)$$

Where:

$$\mathcal{L}(\theta, \phi, X) = \mathcal{E}_{z \sim Q} \left[ \log P_{\theta}(z \mid X) - \log Q_{\phi}(z) \right] \tag{6}$$

#### Evidence lower bound

$$\log P_{\theta}(X^{i}) \ge \mathcal{L}(\theta, \phi, X^{i})$$

$$= \mathcal{E}_{Q_{\phi}(z|X)}[-\log Q_{\phi}(z|X) + \log P_{\theta}(z, X)]$$
(7)

# Relationship between Q and P

The Evidence lower bound (ELBO) or just lower variational bound:

$$\mathcal{D}\left(Q(z) \parallel P(z \mid X)\right) = \mathcal{E}_{z \sim Q}\left[\log Q(z) - \log P(z \mid X)\right] \quad (8)$$

Introduce P(x) and  $P(X \mid z)$ :

$$\mathcal{D}(Q(z) \parallel P(z \mid X)) = \mathcal{E}_{z \sim Q} \left[ \log Q(z) - \log P(z \mid X) - \log P(z) \right] + \log P(X)$$
(9)

Rearrange the ELBO and get another  $\mathcal{KL}$ -term:

$$\log P(X) - \mathcal{D}(Q(z) \parallel P(z \mid X))$$

$$= \mathcal{E}_{z \sim Q} \left[ \log P(X \mid z) \right] - \mathcal{D}(Q(z) \parallel P(z))$$
(10)

And introduce a dependency of Q(z) on X to make it useful:

$$\log P(X) - \mathcal{D}(Q(z \mid X) \parallel P(z \mid X))$$

$$= \mathcal{E}_{z \sim Q} [\log P(X \mid z)] - \mathcal{D}(Q(z \mid X) \parallel P(z))$$
(11)

# Autoencoding variational Bayes

Backpropagation only works with continuous operations, stochastic units aren't that.

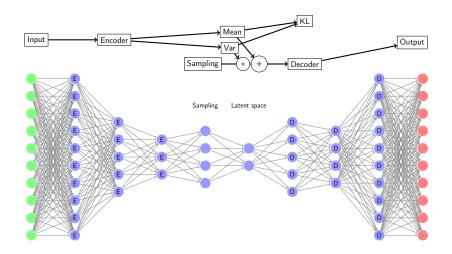
### The reparametrisation trick

With  $\mu(X)$  and  $\sigma(x)$  parametrising  $Q(z \mid X)$  sampling from  $\mathcal{N}(\mu(X), \sigma(X))$  we can draw  $\epsilon \sim \mathcal{N}(0, 1)$  and compute  $z_{\epsilon \sim Q} = \mu(X) + \sqrt{\sigma}(X) \cdot \epsilon$  and perform SGD

$$\mathcal{E}_{z \sim \mathcal{D}} \left[ \mathcal{E}_{\epsilon \sim \mathcal{N}(0,1)} \left[ \log P(X \mid z_{\epsilon \sim Q} - \mathcal{D} \left( Q(z \mid X) \parallel P(z) \right) \right] \right]$$
(12)  
$$\mathcal{L}(\theta, \phi, X) = \mathcal{E}_{z \sim Q} \left[ \log P_{\theta}(z \mid X) - \log Q_{\phi}(z) \right]$$
$$\approx \tilde{\mathcal{L}}^{m}(\theta, \phi, X^{m}) = \frac{n}{m} \sum_{i=1}^{m} \mathcal{L}(\theta, \phi, X_{i})$$
(13)

#### Pseudocode

```
\begin{array}{l} \theta, \phi \leftarrow \text{random parameter} \\ \textbf{repeat} \\ X^m \leftarrow \text{random mini batch with size } m \\ \epsilon \leftarrow \text{random samples from noise distribution } P(\epsilon) \\ g \leftarrow \nabla_{\theta,\phi} \tilde{\mathcal{L}}^m(\theta,\phi,X^m,\epsilon) \\ \theta, \phi \leftarrow \text{update parameters with } g \\ \textbf{until convergence of params } \theta, \phi \\ \textbf{return } \theta, \phi \end{array}
```



## Implementation

- ▶ Using tensorflow.keras functional API [2]
- Currently relu activations and unbiased layers
- Hidden layers constructed iterative
- Deploying a .\_build method, returns encoder, decoder and z
- Model gets trained in customised step
- Sampling layer forms an extra Class

## Training step

```
def train step(self, data):
   with tf.GradientTape() as tape:
       z_mean, z_logvar, z = self.encoder(data)
       reconstruction = self.decoder(z)
       reconstruction_loss = mean(K.square(data - reconstruction),
                                    axis=1)
       kl loss = -.5 * (1 + z logvar - square(z mean) - exp(
           z logvar)
       kl loss = sum(kl loss)
       total loss = kl loss + reconstruction loss
   grads = tape.gradient(total_loss, self.trainable_weights)
   self.optimizer.apply_gradients(zip(grads, self.trainable_weights))
   self.total loss tracker.update state(total loss)
   self.reconstruction_loss_tracker.update_state(reconstruction_loss)
   self.kl_loss_tracker.update_state(kl_loss)
   return {
           "loss": self.total_loss_tracker.result(),
           "reconstruction_loss": self.reconstruction_loss_tracker.result(),
           "kl loss": self.kl loss tracker.result()}
```

# Sampling layer

```
class Sampling_Layer(tf.keras.layers.Layer):
    """
A sampling layer using the mean and log(var) to sample from an input
    """

def call(self, inputs):
    mean, logvar = inputs
    batch = tf.shape(mean)[0]
    dim = tf.shape(mean)[1]
    epsilon = tf.random.normal(shape=(batch, dim), mean=0., stddev=1.)
    return mean + tf.exp(logvar * .5) * epsilon
```

# Training with a dummy

```
import tensorflow as tf
import numpy as np
import generic_VAE
# create some random, uniformly distributed tensors:
training_dummys = np.asarray(
        [tf.random.uniform((400,)) for item in range(100000)])
# create some testing tensors:
testing_dummys = np.asarray(
        [tf.random.uniform((400.)) for item in range(50000)])
# this is completely senseless, the part where a training-set and one for
# testing was created the VAE is unsupervised (in this case)
dummy data = np.concatenate((training dummys, testing dummys))
dummy_vae = generic_VAE.Builder(
        input shape=(400, ).
        encoder_shape=[400, 200, 100, 40, 20, 4],
        decoder_shape=[4, 20, 40, 100, 200, 400],
        latent dims=2.
        dropout rate=0)
vae = generic_VAE.VAE(dummy_vae.decoder_model, dummy_vae.encoder_model)
vae.compile()
vae.fit(dummy_data, epochs=10, batch_size=2000)
```

# Training with a dummy

# Training with GED

[5]

# Next step

[1] []

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  Deep learning with python.

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