

Deep Learning

Lecture 11 – Autoencoders

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Agenda

11.1 Latent Variable Models

11.2 Principal Component Analysis

11.3 Autoencoders

11.4 Variational Autoencoders

11.1

Latent Variable Models

Learning Problems

Supervised Learning:

- ▶ Learn model using dataset with **data-label pairs** $\{(\mathbf{x}_i, \mathbf{y}_i)\}_{i=1}^N$
- ▶ Examples: Classification, regression, structured prediction

Unsupervised Learning:

- ▶ Learn model using dataset **without labels** $\{\mathbf{x}_i\}_{i=1}^N$
- ▶ Examples: Clustering, dimensionality reduction, generative models
- ▶ Some of them use latent variables to capture structure → topic for today

Latent Variable Models

LVMs map between **observation space** $\mathbf{x} \in \mathbb{R}^D$ and **latent space** $\mathbf{z} \in \mathbb{R}^Q$:

$$(f_{\mathbf{w}} : \mathbf{x} \mapsto \mathbf{z}) \quad g_{\mathbf{w}} : \mathbf{z} \mapsto \hat{\mathbf{x}}$$

- ▶ One **latent variable** gets associated with each data point in the training set
- ▶ The latent vectors are smaller than the observations ($Q < D$) \Rightarrow **compression**
- ▶ Models are linear or non-linear, deterministic or stochastic, with/without encoder

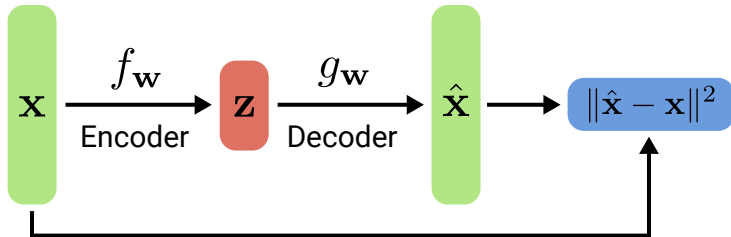
A little taxonomy:

	Deterministic	Probabilistic
Linear	Principle Component Analysis	Probabilistic PCA
Non-Linear w/ Encoder	Autoencoder	Variational Autoencoder
Non-Linear w/o Encoder		Gen. Adv. Networks

Autoencoders

Autoencoders comprise an **encoder** f_w as well as a **decoder** g_w :

$$f_w : \mathbf{x} \mapsto \mathbf{z} \qquad g_w : \mathbf{z} \mapsto \hat{\mathbf{x}}$$

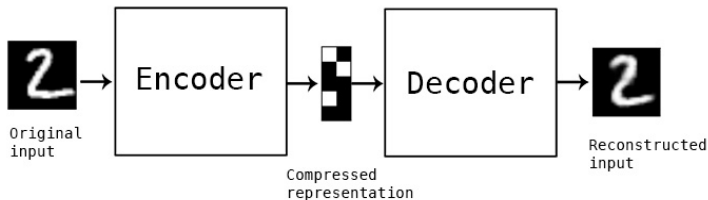


- Models of this type are called **autoencoders** as they predict their input as output
- In contrast, Generative adversarial networks (next lecture) only have a decoder g_w

Autoencoders

Autoencoders comprise an **encoder** $f_{\mathbf{w}}$ as well as a **decoder** $g_{\mathbf{w}}$:

$$f_{\mathbf{w}} : \mathbf{x} \mapsto \mathbf{z} \qquad g_{\mathbf{w}} : \mathbf{z} \mapsto \hat{\mathbf{x}}$$



- ▶ Models of this type are called **autoencoders** as they predict their input as output
- ▶ In contrast, Generative adversarial networks (next lecture) only have a decoder $g_{\mathbf{w}}$

Generative Models

- ▶ Generative modeling is a broad area of machine learning which deals with models of **probability distributions** $p(\mathbf{x})$ over data points \mathbf{x} (e.g., images)
- ▶ The generative model's task is to capture dependencies / structural regularities in the data (e.g., between pixels in images)
- ▶ **Generative latent variable models** capture the structure in **latent variables**
- ▶ Intuitively, we are trying to establish a **theory** for what we observe
- ▶ Some generative models (e.g., normalizing flows) allow for computing $p(\mathbf{x})$
- ▶ Others (e.g., VAEs) only approximate $p(\mathbf{x})$, but allow to draw samples from $p(\mathbf{x})$

Generative Latent Variable Models

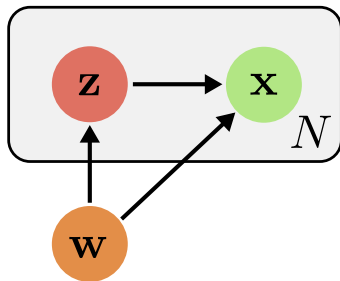
Generative latent variable models often consider a simple **Bayesian model**:

$$p(\mathbf{x}) = \int_{\mathbf{z}} \underbrace{p(\mathbf{z})p(\mathbf{x}|\mathbf{z})}_{\text{Generative Process}} d\mathbf{z} = \mathbb{E}_{\mathbf{z} \sim p(\mathbf{z})} p(\mathbf{x}|\mathbf{z})$$

- ▶ $p(\mathbf{z})$ is the **prior** over the **latent variable** $\mathbf{z} \in \mathbb{R}^Q$
- ▶ $p(\mathbf{x}|\mathbf{z})$ is the **likelihood** (= decoder that transforms \mathbf{z} into a distribution over \mathbf{x})
- ▶ $p(\mathbf{x})$ is the **marginal** of the joint distribution $p(\mathbf{x}, \mathbf{z})$

The goal is to maximize $p(\mathbf{x})$ for a given dataset \mathcal{X} by learning the two models $p(\mathbf{z})$ and $p(\mathbf{x}|\mathbf{z})$ such that the latent variables \mathbf{z} best capture the latent structure of the data.

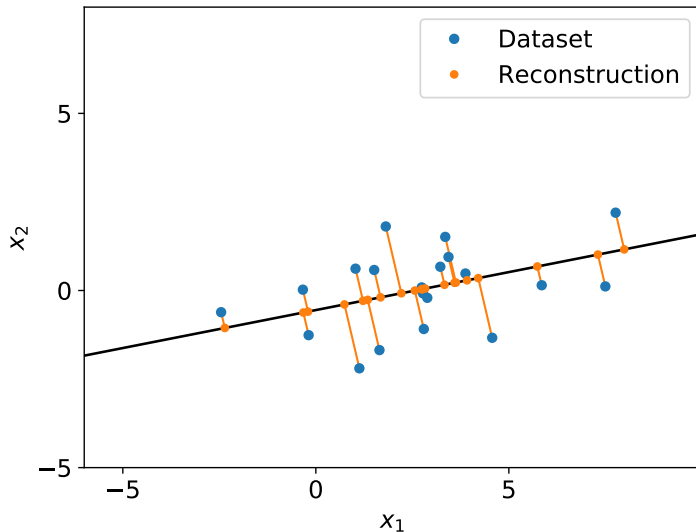
Generative Latent Variable Models



Representation as Graphical Model in Plate Notation:

- ▶ Variables inside plates are replicated (we have N data points to explain)
- ▶ Each data point x is associated with a latent variable z
- ▶ In contrast, parameters are global (exist only once)
- ▶ Remark: We use a single w to refer to all model parameters

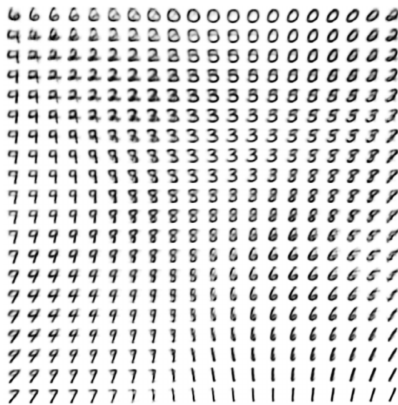
Example: 1D Manifold in 2D Space



Example: Natural Image Manifolds



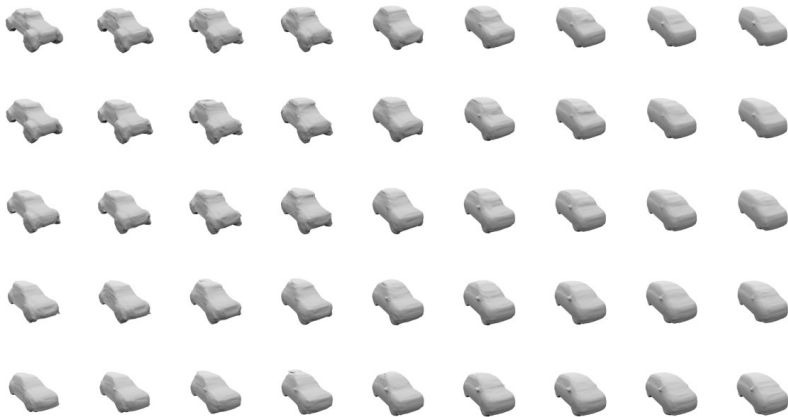
(a) Learned Frey Face manifold



(b) Learned MNIST manifold

- Visualizing the latent space gives insights into the learned semantics

Example: 3D Shape Manifolds



- We can also learn a latent space for 3D shapes and interpolate between them

Example: Sentence Manifolds

“ i want to talk to you . ”

“i want to be with you . ”

“i do n’t want to be with you . ”

i do n’t want to be with you .

she did n’t want to be with him .

he was silent for a long moment .

he was silent for a moment .

it was quiet for a moment .

it was dark and cold .

there was a pause .

it was my turn .

- It is also possible to learn a latent space for sequences of words

11.2

Principal Component Analysis

Preliminaries

- ▶ Let $\mathbf{X} = (\mathbf{x}_1, \dots, \mathbf{x}_N)^\top \in \mathbb{R}^{N \times D}$ denote a **dataset** of observations $\mathbf{x}_i \in \mathbb{R}^D$
- ▶ Let $\mathbf{Z} = (\mathbf{z}_1, \dots, \mathbf{z}_N)^\top \in \mathbb{R}^{N \times Q}$ be the corresponding **latent variables** $\mathbf{z}_i \in \mathbb{R}^Q$
- ▶ While \mathbf{X} is observed, \mathbf{Z} is unobserved and needs to be inferred
- ▶ Typically, we assume $Q < D$, i.e., we want to obtain a **compressed** representation
- ▶ In PCA, our goal is to learn a **linear bidirectional mapping** $\mathcal{Z} \leftrightarrow \mathcal{X}$ such that as much information of \mathcal{X} as possible is retained in \mathcal{Z}
- ▶ In other words, we want to encode $\mathbf{x} \rightarrow \mathbf{z}$ such that if we decode $\mathbf{z} \rightarrow \hat{\mathbf{x}}$, then $\hat{\mathbf{x}}$ is a good approximation to the original \mathbf{x} (in most cases $\hat{\mathbf{x}} \neq \mathbf{x}$)

Principal Component Analysis

Let us assume the following **linear mapping** from latent space to observation space

$$\hat{\mathbf{x}}_i = \bar{\mathbf{x}} + \sum_{j=1}^Q z_{ij} \mathbf{v}_j$$

where $\bar{\mathbf{x}} = \frac{1}{N} \sum_{i=1}^N \mathbf{x}_i$ is the **data mean** and $\mathbf{V} = (\mathbf{v}_1, \dots, \mathbf{v}_Q)$ an **orthonormal basis**.

Our goal is to minimize the L_2 **reconstruction loss** wrt. \mathbf{Z} and \mathbf{V} :

$$\mathcal{L}(\mathbf{Z}, \mathbf{V}) = \sum_{i=1}^N \|\hat{\mathbf{x}}_i - \mathbf{x}_i\|^2 = \sum_{i=1}^N \left\| \underbrace{\bar{\mathbf{x}} + \sum_{j=1}^Q z_{ij} \mathbf{v}_j}_{=\hat{\mathbf{x}}_i} - \mathbf{x}_i \right\|^2$$

Principal Component Analysis

Considering that $\mathbf{V} = (\mathbf{v}_1, \dots, \mathbf{v}_Q)$ is an **orthonormal basis**, we expand \mathcal{L} as follows:

$$\begin{aligned}\mathcal{L}(\mathbf{Z}, \mathbf{V}) &= \sum_{i=1}^N \left\| \bar{\mathbf{x}} + \sum_{j=1}^Q z_{ij} \mathbf{v}_j - \mathbf{x}_i \right\|^2 \\ &= \sum_{i=1}^N \left\| \sum_{j=1}^Q z_{ij} \mathbf{v}_j + \bar{\mathbf{x}} - \mathbf{x}_i \right\|^2 \\ &= \sum_{i=1}^N \left[\sum_{j=1}^Q z_{ij}^2 + 2 \sum_{j=1}^Q z_{ij} \mathbf{v}_j^\top (\bar{\mathbf{x}} - \mathbf{x}_i) + \|\bar{\mathbf{x}} - \mathbf{x}_i\|^2 \right]\end{aligned}$$

Principal Component Analysis

The **reconstruction loss** can therefore be minimized in closed form wrt. \mathbf{Z} :

$$\mathcal{L}(\mathbf{Z}, \mathbf{V}) = \sum_{i=1}^N \left[\sum_{j=1}^Q \left[z_{ij}^2 + 2z_{ij} \mathbf{v}_j^\top (\bar{\mathbf{x}} - \mathbf{x}_i) \right] + \|\bar{\mathbf{x}} - \mathbf{x}_i\|^2 \right]$$
$$\frac{\partial \mathcal{L}(\mathbf{Z}, \mathbf{V})}{\partial z_{ij}} = 2z_{ij} + 2\mathbf{v}_j^\top (\bar{\mathbf{x}} - \mathbf{x}_i) \stackrel{!}{=} 0$$
$$\Rightarrow z_{ij}^* = -\mathbf{v}_j^\top (\bar{\mathbf{x}} - \mathbf{x}_i)$$

For $\mathbf{Z} = \mathbf{Z}^*$, the reconstruction loss simplifies to:

$$\mathcal{L}(\mathbf{Z}^*, \mathbf{V}) = \sum_{i=1}^N \left[- \sum_{j=1}^Q z_{ij}^{*2} + \|\bar{\mathbf{x}} - \mathbf{x}_i\|^2 \right]$$

Principal Component Analysis

The **reconstruction loss** at $z_{ij}^* = -\mathbf{v}_j^\top (\bar{\mathbf{x}} - \mathbf{x}_i)$ can be rewritten as

$$\begin{aligned}\mathcal{L}(\mathbf{Z}^*, \mathbf{V}) &= \sum_{i=1}^N \left[- \sum_{j=1}^Q z_{ij}^{*2} + \|\bar{\mathbf{x}} - \mathbf{x}_i\|^2 \right] \\ &= - \sum_{j=1}^Q \mathbf{v}_j^\top \mathbf{S} \mathbf{v}_j + \sum_{i=1}^N \|\bar{\mathbf{x}} - \mathbf{x}_i\|^2\end{aligned}$$

with \mathbf{S} the **scatter matrix** (unnormalized sample covariance matrix) of \mathbf{x} :

$$\mathbf{S} = \sum_{i=1}^N (\bar{\mathbf{x}} - \mathbf{x}_i)(\bar{\mathbf{x}} - \mathbf{x}_i)^\top$$

Principal Component Analysis

To enforce $\|\mathbf{v}_j\| = 1$, we introduce **Lagrange multipliers** λ_j into the loss:

$$\begin{aligned}\mathcal{L}(\mathbf{Z}^*, \mathbf{V}, \boldsymbol{\lambda}) &= -\sum_{j=1}^Q \mathbf{v}_j^\top \mathbf{S} \mathbf{v}_j + \sum_{i=1}^N \|\bar{\mathbf{x}} - \mathbf{x}_i\|^2 + \sum_{j=1}^Q \lambda_j (\mathbf{v}_j^\top \mathbf{v}_j - 1) \\ \frac{\partial \mathcal{L}(\mathbf{Z}^*, \mathbf{V}, \boldsymbol{\lambda})}{\partial \mathbf{v}_j} &= -2\mathbf{S} \mathbf{v}_j + 2\lambda_j \mathbf{v}_j \stackrel{!}{=} 0 \\ \Rightarrow \mathbf{S} \mathbf{v}_j &= \lambda_j \mathbf{v}_j\end{aligned}$$

We see that $\{\boldsymbol{\lambda}, \mathbf{V}\}$ is the solution to an **eigenvalue problem**.

We also observe that as we have $\mathbf{v}_j^\top \mathbf{S} \mathbf{v}_j = \lambda_j \mathbf{v}_j^\top \mathbf{v}_j = \lambda_j$, the loss \mathcal{L} is minimized by choosing (for \mathbf{V}) the eigenvectors \mathbf{v}_j of \mathbf{S} corresponding to the top Q eigenvalues.

Principal Component Analysis

Consider again the **linear model** that we started with:

$$\hat{\mathbf{x}}_i = \bar{\mathbf{x}} + \sum_{j=1}^Q z_{ij} \mathbf{v}_j$$

Both the PCA decoder and encoder are simple **linear mappings**:

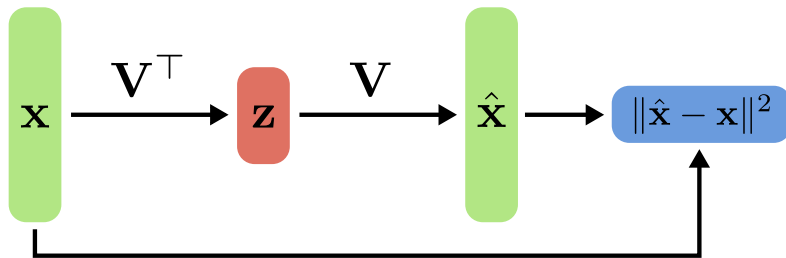
Decoder:

$$\mathbf{x} = \mathbf{V}\mathbf{z} + \bar{\mathbf{x}}$$

Encoder:

$$\mathbf{z} = \mathbf{V}^\top (\mathbf{x} - \bar{\mathbf{x}})$$

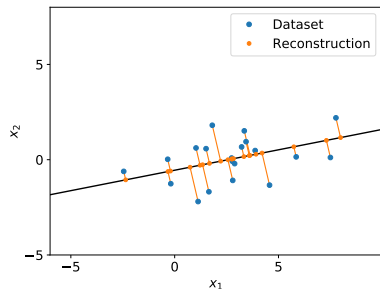
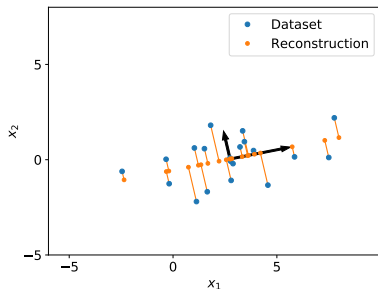
Principal Component Analysis



PCA Recipe:

- ▶ Given a dataset $\mathcal{X} = \{\mathbf{x}_1, \dots, \mathbf{x}_N\}$ of observations $\mathbf{x}_i \in \mathbb{R}^D$
- ▶ Compute the **data mean** $\bar{\mathbf{x}}$ and **scatter matrix** $\mathbf{S} = \sum_{i=1}^N (\bar{\mathbf{x}} - \mathbf{x}_i)(\bar{\mathbf{x}} - \mathbf{x}_i)^\top$
- ▶ Compute the **eigen decomposition** of \mathbf{S}
- ▶ Select the Q eigenvectors corresponding to the Q largest eigenvalues for \mathbf{V}

Principal Component Analysis



There are 2 perspectives on PCA:

- ▶ We saw that PCA can be motivated by **minimizing the L_2 reconstruction error**
- ▶ However, we can also motivate PCA by **maximizing the variance** of latent points
- ▶ In other words, we like to find an embedding that captures most of the variation in the original dataset while using a smaller dimensionality $Q \ll D$

Variance Maximization Perspective

Consider the following (one-dimensional) encoding of vector \mathbf{x} :

$$\mathbf{z} = \mathbf{v}^\top (\mathbf{x} - \bar{\mathbf{x}})$$

Our goal is to **maximize variance** in latent space:

$$\begin{aligned}\text{Var}(\mathbf{z}) &= \mathbb{E} \left[\left(\mathbf{v}^\top (\mathbf{x} - \bar{\mathbf{x}}) - \mathbb{E} \left[\mathbf{v}^\top (\mathbf{x} - \bar{\mathbf{x}}) \right] \right)^2 \right] \\ &= \mathbb{E} \left[\left(\mathbf{v}^\top (\mathbf{x} - \bar{\mathbf{x}}) \right)^2 \right] \quad (\text{as } \mathbb{E}[\mathbf{x}] = \bar{\mathbf{x}}) \\ &= \mathbb{E} \left[\mathbf{v}^\top (\bar{\mathbf{x}} - \mathbf{x})(\bar{\mathbf{x}} - \mathbf{x})^\top \mathbf{v} \right] \\ &\propto \mathbf{v}^\top \mathbf{S} \mathbf{v} \quad (\text{as } \mathbf{S} \text{ is not normalized})\end{aligned}$$

Variance Maximization Perspective

Let us now again assume an orthonormal basis $\mathbf{V} = (\mathbf{v}_1, \dots, \mathbf{v}_Q)$ of dimension Q .

Maximizing the sum of variances along each dimension subject to normalization constraints leads to the optimization problem we are already familiar with:

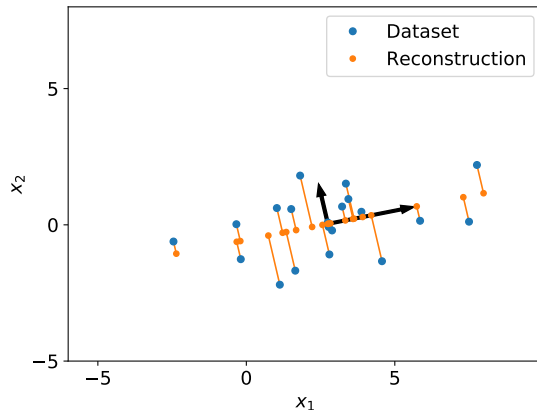
$$\boldsymbol{\lambda}^*, \mathbf{V}^* = \operatorname{argmax}_{\boldsymbol{\lambda}, \mathbf{V}} \sum_{j=1}^Q \mathbf{v}_j^\top \mathbf{S} \mathbf{v}_j + \sum_{j=1}^Q \lambda_j (\mathbf{v}_j^\top \mathbf{v}_j - 1)$$

A solution is given by the Q largest eigenvalues and corresponding eigenvectors of \mathbf{S} .

Remark: $\mathbf{v}_j^\top \mathbf{S} \mathbf{v}_j = \lambda_j \mathbf{v}_j^\top \mathbf{v}_j = \lambda_j$ is the **variance** along the j 'th principal component if the scatter matrix \mathbf{S} is normalized by the number of data points (=covariance matrix).

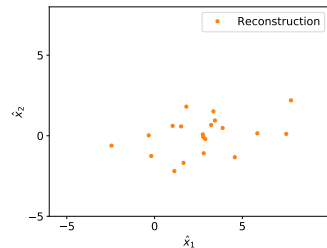
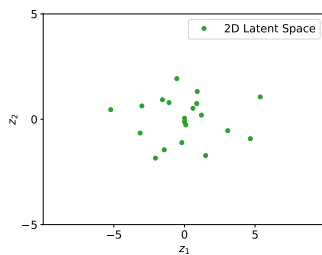
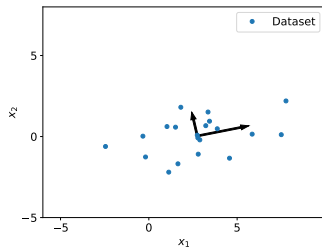
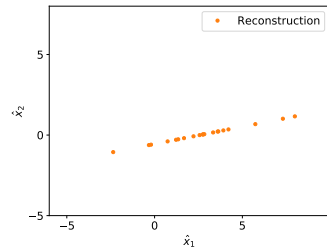
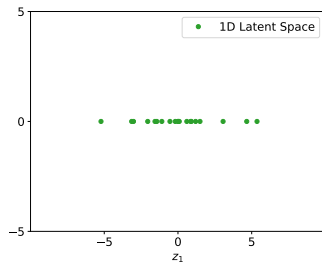
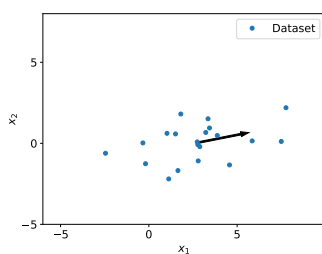
Examples

Results on Synthetic 2D Dataset

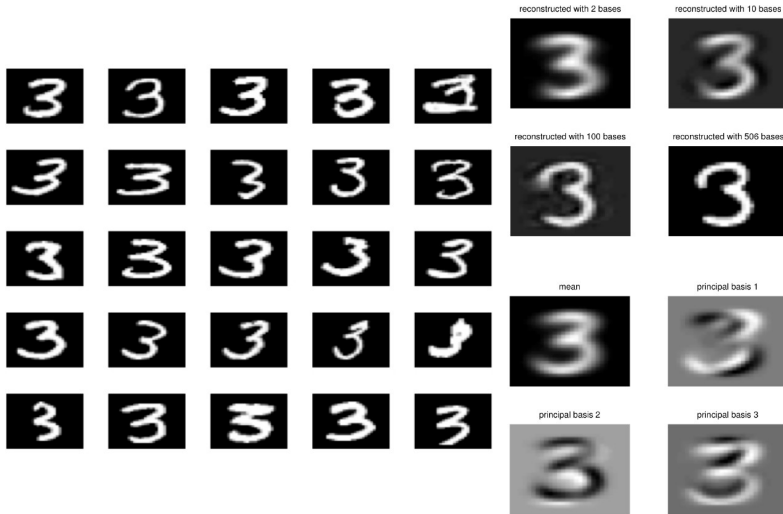


- PCA on a dataset with $N = 20$, $D = 2$ and $Q = 1$ (projection onto 1D subspace)
- The two eigenvectors \mathbf{v}_j shown in black are scaled by $\sqrt{\lambda_j}$

Results on Synthetic 2D Dataset



Results on MNIST



Results on Faces

PCA on Face Images:

- ▶ PCA on 2429 19x19 grayscale images (CBCL data)
- ▶ Yields good reconstructions with only 3 components:



- ▶ We can apply a classifier directly on this latent representation
- ▶ PCA with 3 components obtains 79 % accuracy on face/non-face discrimination on test data vs. 76.8 % for a Gaussian Mixture Model with 84 states
- ▶ Also commonly used for analyzing the latent properties of a dataset

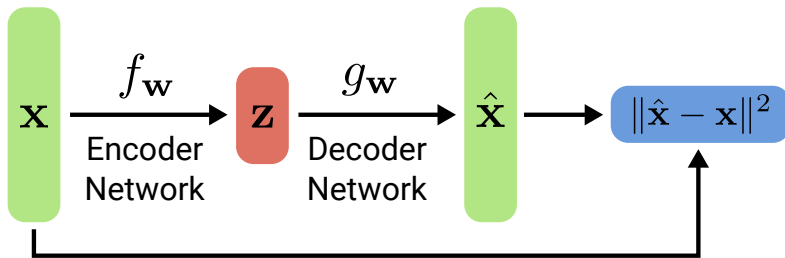
Learned Basis (Eigenfaces)



11.3

Autoencoders

Autoencoders



- ▶ An autoencoder is a **neural network** whose outputs are its own inputs
- ▶ The input $\mathbf{x} \in \mathbb{R}^D$ is compressed to a latent code $\mathbf{z} \in \mathbb{R}^Q$
- ▶ The goal is to **minimize the reconstruction error** (as in PCA)

PCA as Special Case of Autoencoders

Let $f_{\mathbf{w}} : \mathbf{x} \mapsto \mathbf{z}$ denote the encoder and $g_{\mathbf{w}} : \mathbf{z} \mapsto \hat{\mathbf{x}}$ the decoder network.
Let's assume both mappings to be **linear** (without activation function):

$$\mathbf{z} = f_{\mathbf{w}}(\mathbf{x}) = \mathbf{A}\mathbf{x} + \mathbf{a} \quad \hat{\mathbf{x}} = g_{\mathbf{w}}(\mathbf{z}) = \mathbf{B}\mathbf{z} + \mathbf{b}$$

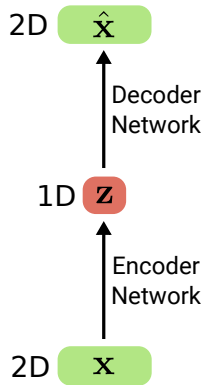
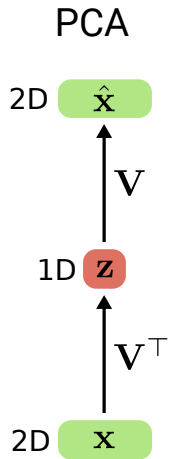
In this case, we have:

$$\hat{\mathbf{x}} = g_{\mathbf{w}}(f_{\mathbf{w}}(\mathbf{x})) = \underbrace{\mathbf{B}\mathbf{A}}_{=\mathbf{C}}\mathbf{x} + \underbrace{\mathbf{a} + \mathbf{b}}_{=\mathbf{c}}$$

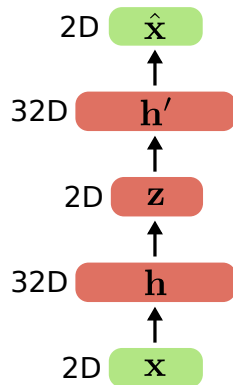
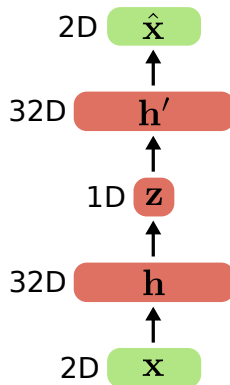
Thus, the optimal solution \mathbf{w}^* is given by **PCA**:

$$\mathbf{w}^* = \underset{\mathbf{w}}{\operatorname{argmin}} \sum_{i=1}^N \left\| \underbrace{\mathbf{C}\mathbf{x}_i + \mathbf{c}}_{=\hat{\mathbf{x}}_i} - \mathbf{x}_i \right\|^2$$

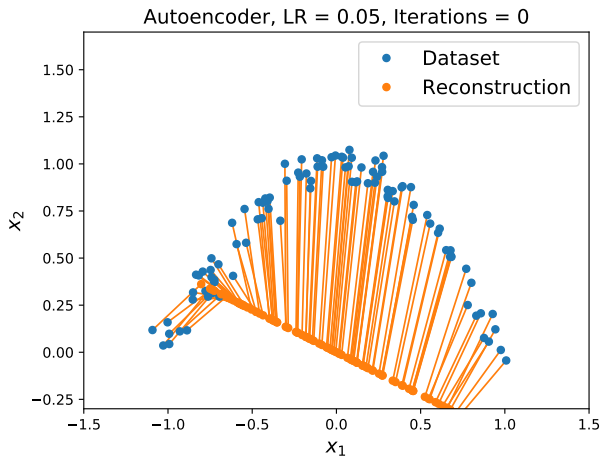
Results on Synthetic 2D Dataset



Autoencoders

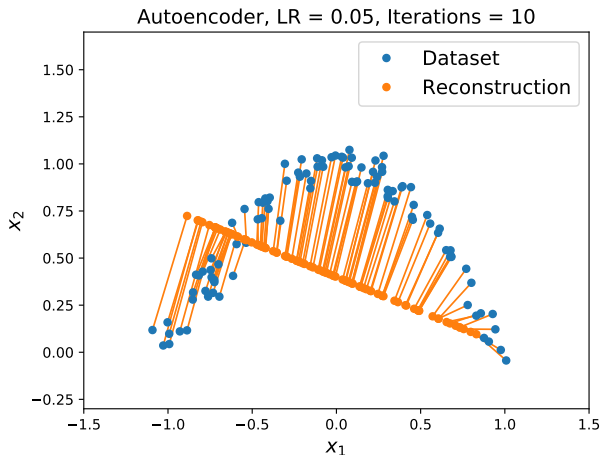


Results on Synthetic 2D Dataset



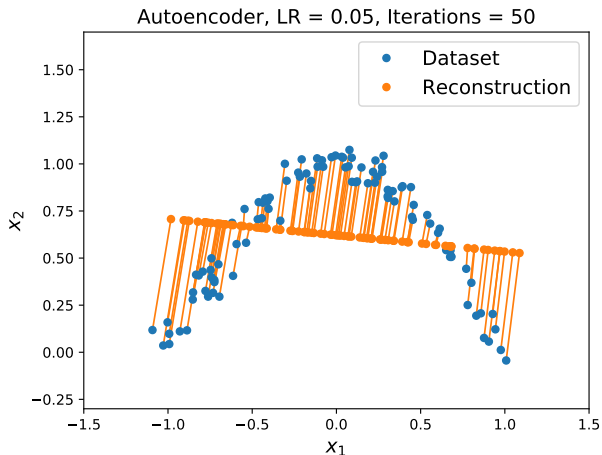
- **Autoencoder reconstructions** using one linear encoder/decoder layer with $Q = 1$

Results on Synthetic 2D Dataset



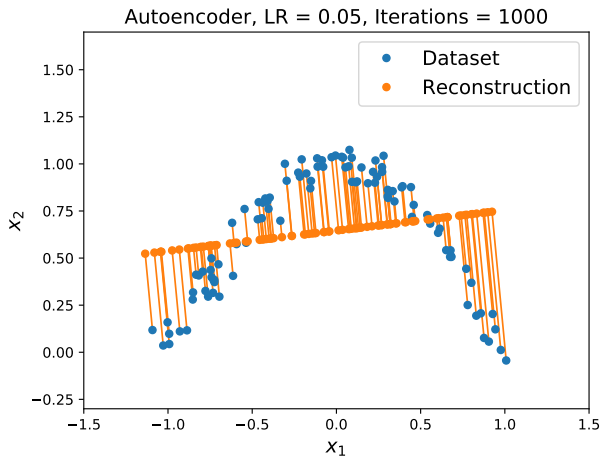
► **Autoencoder reconstructions** using one linear encoder/decoder layer with $Q = 1$

Results on Synthetic 2D Dataset



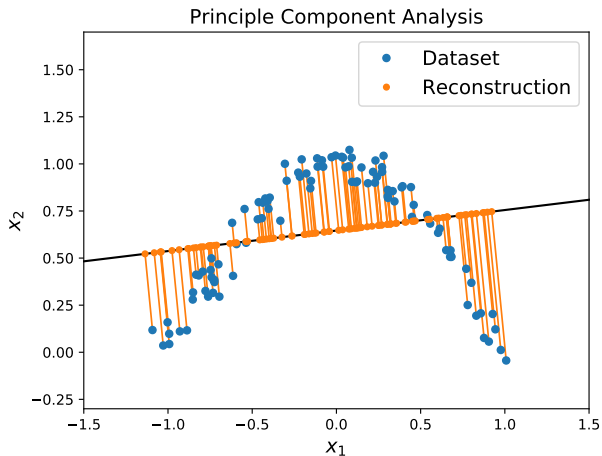
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Results on Synthetic 2D Dataset



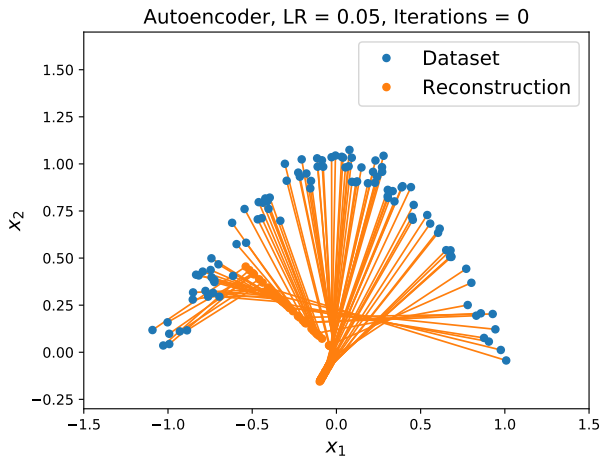
- **Autoencoder reconstructions** using one linear encoder/decoder layer with $Q = 1$

Results on Synthetic 2D Dataset



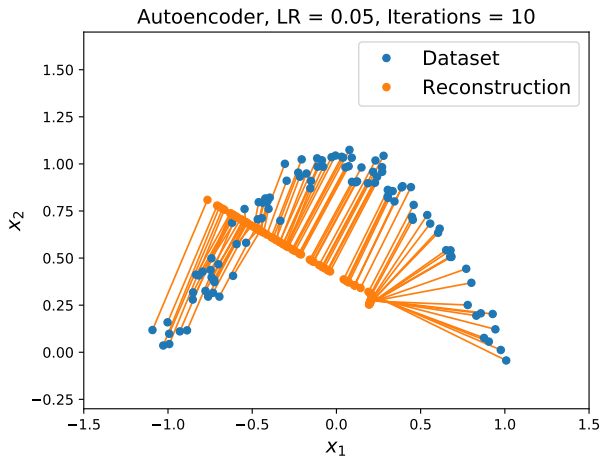
► **PCA reconstructions** on the same dataset with $Q = 1$

Results on Synthetic 2D Dataset



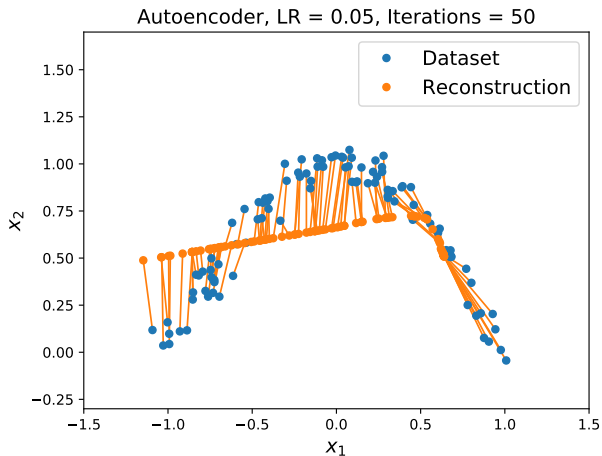
- **Autoencoder reconstructions** using 32 dimensional hidden layers and $Q = 1$

Results on Synthetic 2D Dataset



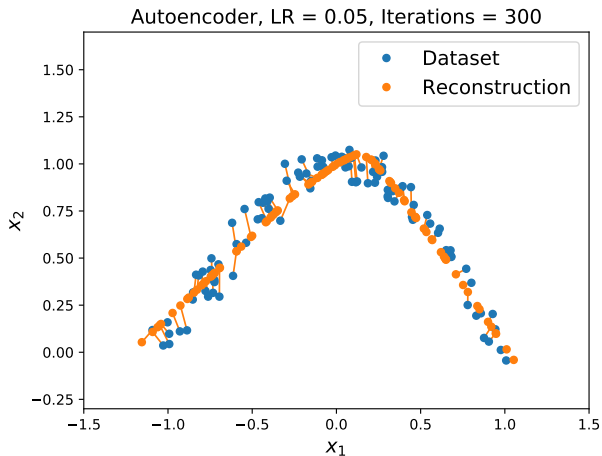
- **Autoencoder reconstructions** using 32 dimensional hidden layers and $Q = 1$

Results on Synthetic 2D Dataset



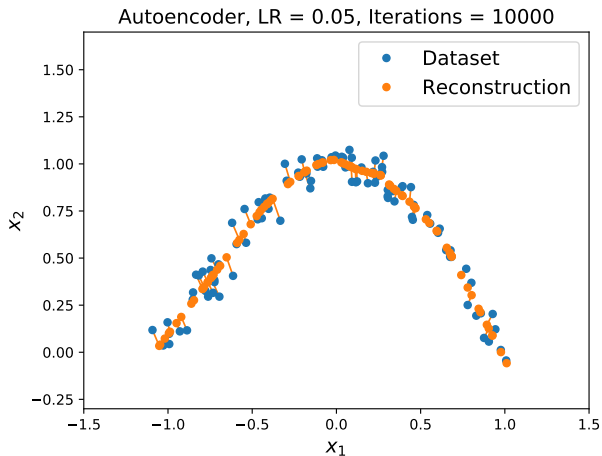
- **Autoencoder reconstructions** using 32 dimensional hidden layers and $Q = 1$

Results on Synthetic 2D Dataset



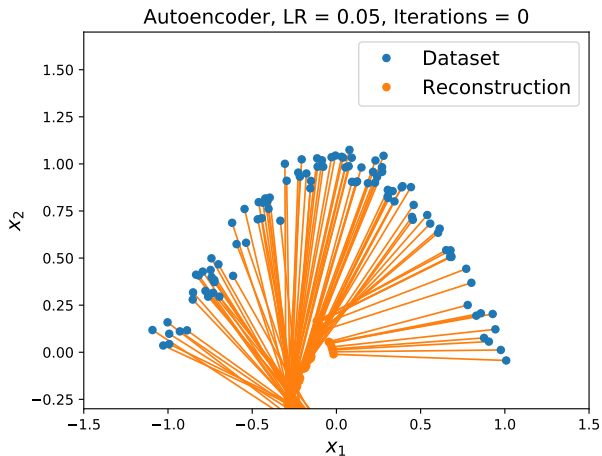
- **Autoencoder reconstructions** using 32 dimensional hidden layers and $Q = 1$

Results on Synthetic 2D Dataset



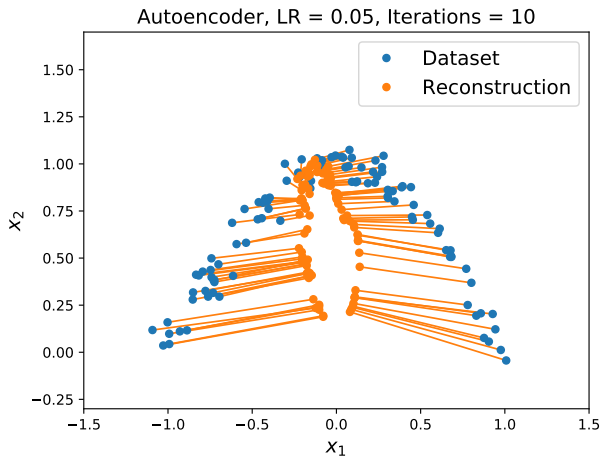
- **Autoencoder reconstructions** using 32 dimensional hidden layers and $Q = 1$

Results on Synthetic 2D Dataset



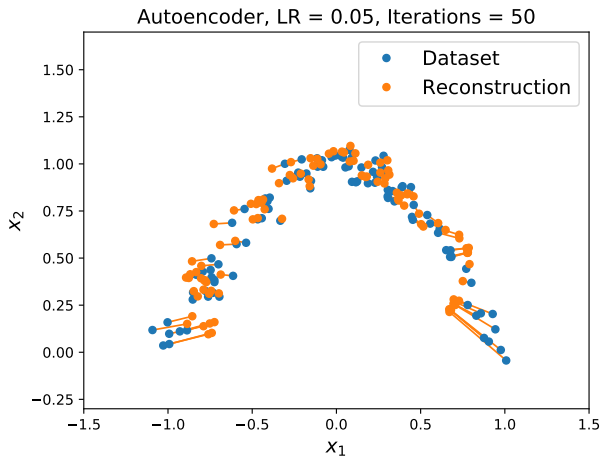
- **Autoencoder reconstructions** using 32 dimensional hidden layers and $Q = 2$

Results on Synthetic 2D Dataset



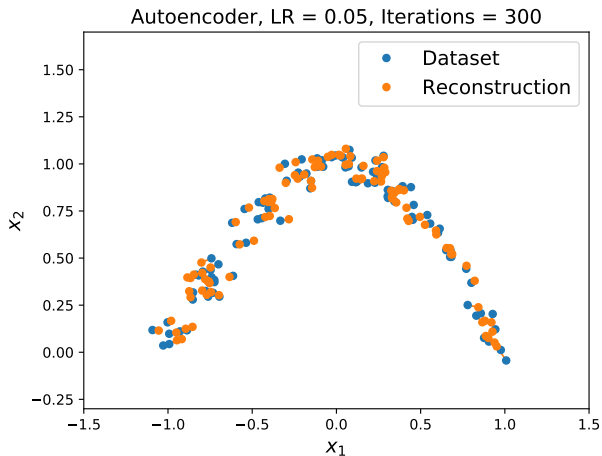
- **Autoencoder reconstructions** using 32 dimensional hidden layers and $Q = 2$

Results on Synthetic 2D Dataset



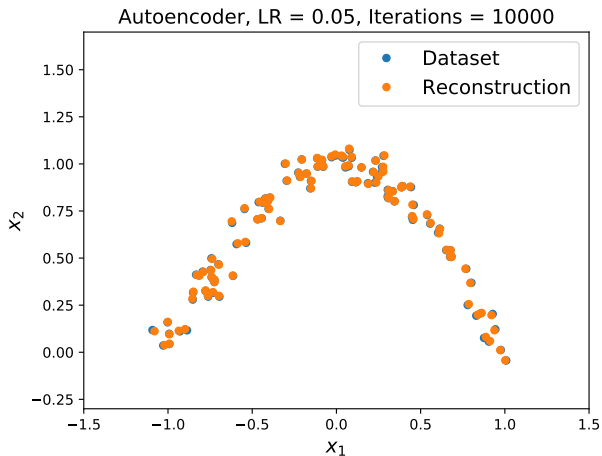
- **Autoencoder reconstructions** using 32 dimensional hidden layers and $Q = 2$

Results on Synthetic 2D Dataset



- **Autoencoder reconstructions** using 32 dimensional hidden layers and $Q = 2$

Results on Synthetic 2D Dataset



- **Autoencoder reconstructions** using 32 dimensional hidden layers and $Q = 2$

Comparing Reconstructions



Real data

30-d deep autoencoder

30-d logistic PCA

30-d PCA

- ▶ Deep autoencoders learn **non-linear** latent embeddings
- ▶ They often have **smaller reconstruction errors** compared to PCA with same Q
- ▶ In contrast, PCA always learns the best linear mapping

Denoising Autoencoders



- ▶ **Denoising Autoencoders** take noisy inputs and predict the original noise-free data
- ▶ Higher level representations are relatively stable and robust to input corruption
- ▶ Encourages the model to **generalize better** and capture useful structure
- ▶ Similar to data augmentation (except that the “label” is the noise-free input)
- ▶ <https://blog.keras.io/building-autoencoders-in-keras.html>

11.4

Variational Autoencoders

A Bayesian Generative Latent Variable Model

So far, we have discussed deterministic latent variables. We will now take a **probabilistic perspective** on **latent variable models** with autoencoding properties. Consider the following **Bayesian model** of the data $\mathbf{x} \in \mathbb{R}^D$:

$$p(\mathbf{x}) = \int_{\mathbf{z}} \underbrace{p(\mathbf{z})p(\mathbf{x}|\mathbf{z})}_{\text{Generative Process}} d\mathbf{z} = \mathbb{E}_{\mathbf{z} \sim p(\mathbf{z})} p(\mathbf{x}|\mathbf{z})$$

- ▶ $p(\mathbf{z})$ is the **prior** over the **latent variable** $\mathbf{z} \in \mathbb{R}^Q$
- ▶ $p(\mathbf{x}|\mathbf{z})$ is the **likelihood**
- ▶ $p(\mathbf{x})$ is the **marginal** of the joint distribution $p(\mathbf{x}, \mathbf{z})$

A Bayesian Generative Latent Variable Model

Assumptions:

- ▶ We assume the **prior** model $p(\mathbf{z})$ to be samplable and computable
- ▶ We assume the **likelihood** model $p(\mathbf{x}|\mathbf{z})$ to be computable
- ▶ In other words, we can **sample** from $p(\mathbf{z})$ and we can **compute** the probability mass/density of $p(\mathbf{z})$ and $p(\mathbf{x}|\mathbf{z})$ for any given \mathbf{x} and \mathbf{z}
- ▶ These assumptions hold for autoregressive models (e.g., language models)
- ▶ However, they fail for loopy graphical models where approximations must be used
- ▶ We will choose **simple parameteric distributions** to achieve this

A Bayesian Generative Latent Variable Model

To find the model parameters \mathbf{w} , we want to minimize the **negative log likelihood**:

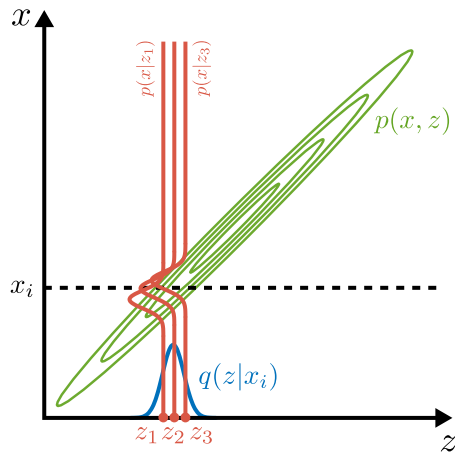
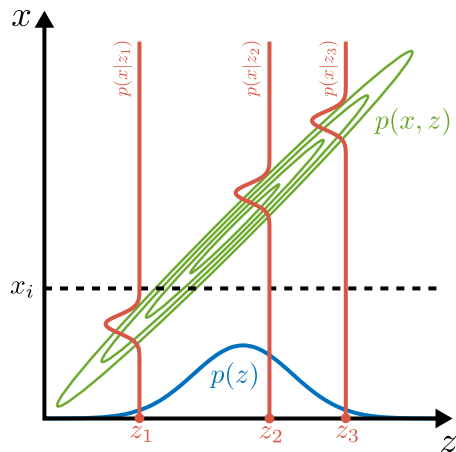
$$\begin{aligned}\mathbf{w}^* &= \operatorname{argmin}_{\mathbf{w}} \mathbb{E}_{\mathbf{x} \sim p_{data}} [-\log p_{\mathbf{w}}(\mathbf{x})] \\ &= \operatorname{argmin}_{\mathbf{w}} \mathbb{E}_{\mathbf{x} \sim p_{data}} [-\log \mathbb{E}_{\mathbf{z} \sim p_{\mathbf{w}}(\mathbf{z})} p_{\mathbf{w}}(\mathbf{x}|\mathbf{z})] \\ &= \operatorname{argmin}_{\mathbf{w}} \sum_{i=1}^N -\log \mathbb{E}_{\mathbf{z} \sim p_{\mathbf{w}}(\mathbf{z})} p_{\mathbf{w}}(\mathbf{x}_i|\mathbf{z})\end{aligned}$$

- ▶ Unfortunately, even given our assumption, computing $p_{\mathbf{w}}(\mathbf{x})$ is **intractable**
- ▶ VAEs side-step this by introducing another model component, a so-called **recognition model** $q_{\mathbf{w}}(\mathbf{z}|\mathbf{x})$ to approximate the true posterior $p_{\mathbf{w}}(\mathbf{z}|\mathbf{x})$

Intuition for Intractability

- ▶ Imagine the observations are sound waves and the latents are word sequences
- ▶ We are looking for a “theory” of sound waves that best explains them
- ▶ We want to minimize the cross entropy between the data distribution over sound waves and our model distribution of the sound waves
- ▶ However, the marginal $p_{\mathbf{w}}(\mathbf{x})$ is intractable due to the large search space
- ▶ If we listen to a song, it is sometimes unclear what the lyrics are
- ▶ If someone tells us the lyrics, we can suddenly hear it / verify it
- ▶ But the search over the word sequences that explain the sound waves is hard as there are many sequences and we might not think of the right one
- ▶ VAEs thus use a recognition model $q_{\mathbf{w}}(\mathbf{z}|\mathbf{x})$ that computes the word sequence
- ▶ This model does not need to be correct, it is an approximation to $p(\mathbf{z}|\mathbf{x})$

Intuition for Intractability



- Computing $p_{\mathbf{w}}(\mathbf{x}) = \mathbb{E}_{\mathbf{z}} p_{\mathbf{w}}(\mathbf{x}|\mathbf{z})$ is hard, in particular in high dimensions

The Evidence Lower Bound (ELBO)

We seek a **tractable lower bound** to the likelihood:

$$\begin{aligned}\log p(\mathbf{x}) &= \mathbb{E}_{\mathbf{z} \sim q(\mathbf{z}|\mathbf{x})} \log \frac{p(\mathbf{x})p(\mathbf{z}|\mathbf{x})}{p(\mathbf{z}|\mathbf{x})} \\ &= \mathbb{E}_{\mathbf{z} \sim q(\mathbf{z}|\mathbf{x})} \log \frac{p(\mathbf{z}, \mathbf{x})}{q(\mathbf{z}|\mathbf{x})} + \log \frac{q(\mathbf{z}|\mathbf{x})}{p(\mathbf{z}|\mathbf{x})} \\ &= \mathbb{E}_{\mathbf{z} \sim q(\mathbf{z}|\mathbf{x})} \log \frac{p(\mathbf{z}, \mathbf{x})}{q(\mathbf{z}|\mathbf{x})} + \underbrace{KL(q(\mathbf{z}|\mathbf{x}), p(\mathbf{z}|\mathbf{x}))}_{\geq 0} \\ &\geq \mathbb{E}_{\mathbf{z} \sim q(\mathbf{z}|\mathbf{x})} \log \frac{p(\mathbf{z}, \mathbf{x})}{q(\mathbf{z}|\mathbf{x})} \quad (\text{ELBO})\end{aligned}$$

- ▶ In practice, $q(\mathbf{z}|\mathbf{x})$ is a **variational approximation** to the true posterior $p(\mathbf{z}|\mathbf{x})$
- ▶ Therefore, the ELBO is sometimes also called **variational lower bound**
- ▶ The divergence term $KL(q(\mathbf{z}|\mathbf{x}), p(\mathbf{z}|\mathbf{x}))$ measures the **approximation error**

The Evidence Lower Bound (ELBO)

The **negative log likelihood** is thus **upper bounded** by:

$$\begin{aligned} -\log p(\mathbf{x}) &\leq \mathbb{E}_{\mathbf{z} \sim q(\mathbf{z}|\mathbf{x})} \log \frac{q(\mathbf{z}|\mathbf{x})}{p(\mathbf{z}, \mathbf{x})} \\ &= \mathbb{E}_{\mathbf{z} \sim q(\mathbf{z}|\mathbf{x})} \log \frac{q(\mathbf{z}|\mathbf{x})}{p(\mathbf{z})p(\mathbf{x}|\mathbf{z})} \\ &= \mathbb{E}_{\mathbf{z} \sim q(\mathbf{z}|\mathbf{x})} \log \frac{q(\mathbf{z}|\mathbf{x})}{p(\mathbf{z})} - \log p(\mathbf{x}|\mathbf{z}) \\ &= KL(q(\mathbf{z}|\mathbf{x}), p(\mathbf{z})) + \mathbb{E}_{\mathbf{z} \sim q(\mathbf{z}|\mathbf{x})} [-\log p(\mathbf{x}|\mathbf{z})] \end{aligned}$$

- The bound comprises a **KL divergence** and a **conditional log likelihood**
- Note how $q(\mathbf{z}|\mathbf{x})$ and $p(\mathbf{x}|\mathbf{z})$ act as an **autoencoder** model: $\mathbf{x} \xrightarrow{q} \mathbf{z} \xrightarrow{p} \mathbf{x}$

Variational Autoencoder (VAE)

Let $\mathcal{X} = \{\mathbf{x}_1, \dots, \mathbf{x}_N\}_{i=1}^N$ with $\mathbf{x}_i \in \mathbb{R}^D$ be a dataset and \mathbf{w} the model parameters.

The **Variational Autoencoder** minimizes this bound to the **negative log likelihood**:

$$\begin{aligned}\mathbf{w}^* &= \operatorname{argmin}_{\mathbf{w}} \mathbb{E}_{\mathbf{x} \sim p_{data}} [-\log p_{\mathbf{w}}(\mathbf{x})] \\ &= \operatorname{argmin}_{\mathbf{w}} \sum_{i=1}^N [-\log p_{\mathbf{w}}(\mathbf{x}_i)] \\ &\approx \operatorname{argmin}_{\mathbf{w}} \underbrace{\sum_{i=1}^N KL(q_{\mathbf{w}}(\mathbf{z}|\mathbf{x}_i), p(\mathbf{z}))}_{\text{Approx. Posterior = Prior}} + \underbrace{\mathbb{E}_{\mathbf{z} \sim q_{\mathbf{w}}(\mathbf{z}|\mathbf{x}_i)} [-\log p_{\mathbf{w}}(\mathbf{x}_i|\mathbf{z})]}_{\text{Reconstruction Term}}\end{aligned}$$

- ▶ In a VAE, $q_{\mathbf{w}}(\mathbf{z}|\mathbf{x})$ is a **multivariate Gaussian** parameterized by a neural network
- ▶ It thus makes a **variational approximation** $q_{\mathbf{w}}(\mathbf{z}|\mathbf{x})$ to the true posterior $p(\mathbf{z}|\mathbf{x})$

Variational Autoencoder (VAE)

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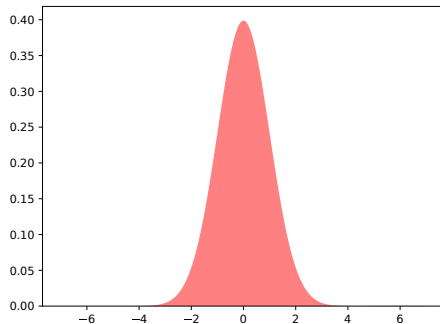
- The **likelihood model** $p_{\mathbf{w}}(\mathbf{x}_i|\mathbf{z})$ is also parameterized by a neural network
- For binary observations **Bernoulli**, for real observations **Gaussian** or **Laplacian**

Neural Network Parameterization

Let us consider a **Gaussian recognition model**:

$$q_{\mathbf{w}}(\mathbf{z}|\mathbf{x}) = \frac{1}{(2\pi)^{Q/2}} \frac{1}{|\Sigma_{\mathbf{w}}(\mathbf{x})|^{1/2}} \exp \left(-\frac{1}{2} (\mathbf{z} - \mu_{\mathbf{w}}(\mathbf{x}))^\top \Sigma_{\mathbf{w}}(\mathbf{x})^{-1} (\mathbf{z} - \mu_{\mathbf{w}}(\mathbf{x})) \right)$$

- ▶ The mean μ and covariance Σ are functions of \mathbf{x} and with parameters \mathbf{w}
- ▶ In a VAE, these functions are implemented using a **neural network** (e.g., MLP)
- ▶ They often have a **shared backbone**
- ▶ Typically, we use $\Sigma_{\mathbf{w}}(\mathbf{x}) = \text{diag}(\sigma_{\mathbf{w}}^2(\mathbf{x}))$



Recognition Model and Prior

Assume a **Gaussian recognition model** $q(\mathbf{z}) = \mathcal{N}(\mathbf{z}|\boldsymbol{\mu}, \boldsymbol{\sigma}^2)$ and **prior** $p(\mathbf{z}) = \mathcal{N}(\mathbf{0}, \mathbf{I})$.
(note that for clarity, we drop the dependency of q , $\boldsymbol{\mu}$ and $\boldsymbol{\sigma}$ on \mathbf{x} and \mathbf{w})

$$\int q(\mathbf{z}) \log q(\mathbf{z}) d\mathbf{z} = \int \mathcal{N}(\mathbf{z}|\boldsymbol{\mu}, \boldsymbol{\sigma}^2) \log \mathcal{N}(\mathbf{z}|\boldsymbol{\mu}, \boldsymbol{\sigma}^2) d\mathbf{z} = -\frac{Q}{2} \log(2\pi) - \frac{1}{2} \sum_{j=1}^J (1 + \log \sigma_j^2)$$

$$\int q(\mathbf{z}) \log p(\mathbf{z}) d\mathbf{z} = \int \mathcal{N}(\mathbf{z}|\boldsymbol{\mu}, \boldsymbol{\sigma}^2) \log \mathcal{N}(\mathbf{z}|\mathbf{0}, \mathbf{I}) d\mathbf{z} = -\frac{Q}{2} \log(2\pi) - \frac{1}{2} \sum_{j=1}^J (\mu_j^2 + \sigma_j^2)$$

$$KL(q(\mathbf{z}), p(\mathbf{z})) = \int q(\mathbf{z}) (\log q(\mathbf{z}) - \log p(\mathbf{z})) d\mathbf{z} = \frac{1}{2} \sum_{j=1}^J (\mu_j^2 + \sigma_j^2 - 1 - \log \sigma_j^2)$$

The KL term has a simple **analytical solution** in this case. This is the standard setup.
The **covariance matrix** $\boldsymbol{\Sigma} = \text{diag}(\boldsymbol{\sigma}^2)$ of the recognition model is chosen **diagonal**.

Learning a VAE

Variational Autoencoder Objective:

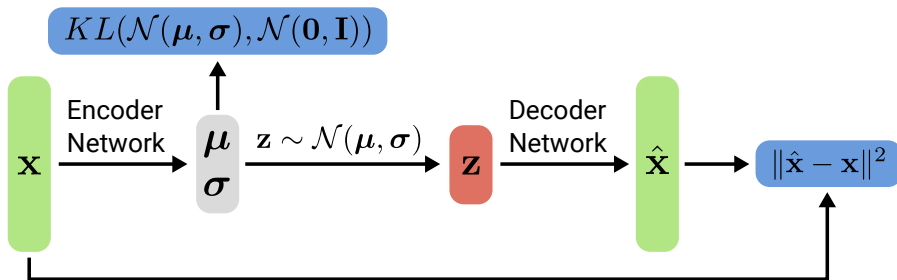
$$\mathbf{w}^* = \underset{\mathbf{w}}{\operatorname{argmin}} \sum_{i=1}^N \underbrace{KL(q_{\mathbf{w}}(\mathbf{z}|\mathbf{x}_i), p(\mathbf{z}))}_{\text{Approx. Posterior = Prior}} + \underbrace{\mathbb{E}_{\mathbf{z} \sim q_{\mathbf{w}}(\mathbf{z}|\mathbf{x}_i)} [-\log p_{\mathbf{w}}(\mathbf{x}_i|\mathbf{z})]}_{\text{Reconstruction Term}}$$

- ▶ The gradients for the **KL term** wrt. \mathbf{w} are easily obtained using backpropagation
- ▶ For the **reconstruction term**, the forward pass can be computed using sampling, but the backward pass requires differentiating through a sampling operation
- ▶ Solved by the **reparameterization trick** which moves the sampling to the input:

$$\mathbb{E}_{\mathbf{z} \sim q_{\mathbf{w}}(\mathbf{z}|\mathbf{x}_i)} [-\log p_{\mathbf{w}}(\mathbf{x}_i|\mathbf{z})] = \mathbb{E}_{\boldsymbol{\epsilon} \sim \mathcal{N}(\mathbf{0}, \mathbf{I})} [-\log p_{\mathbf{w}}(\mathbf{x}_i|\mathbf{z} = \boldsymbol{\mu}_{\mathbf{w}}(\mathbf{x}_i) + \boldsymbol{\sigma}_{\mathbf{w}}(\mathbf{x}_i) \odot \boldsymbol{\epsilon})]$$

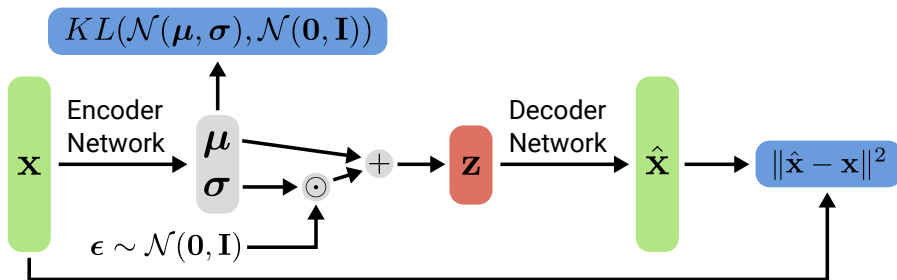
- ▶ In practice, a single sample $\boldsymbol{\epsilon}$ per \mathbf{x} often suffices (depends on minibatch size)

Learning a VAE



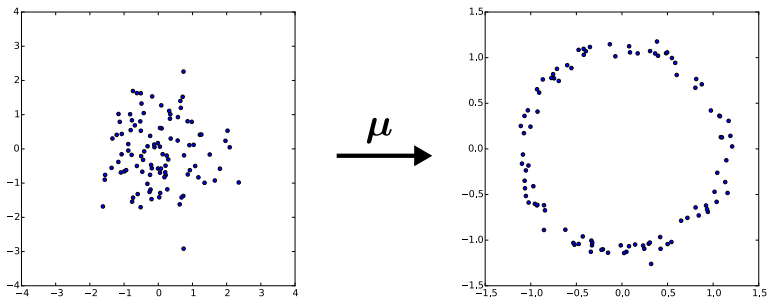
- **Vanilla VAE** formulation which is intractable due to sampling inside the network
- Remark: We assume a Gaussian likelihood $p_{\mathbf{w}}(\mathbf{x}|\mathbf{z})$ in this example

Learning a VAE



- **Reparameterized version** which is tractable as sampling has been moved to input
- This trick works for Gaussians and some other simple distributions (cf., Kingma)

Expressiveness

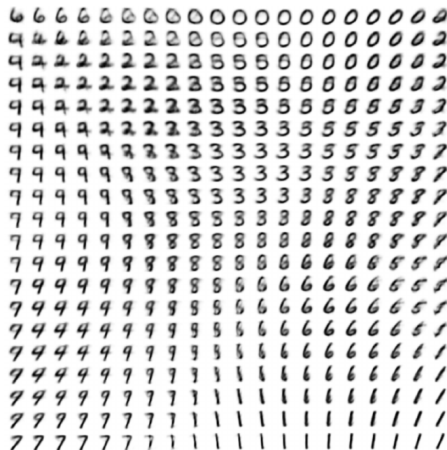


- ▶ VAEs are very **expressive**: Consider random samples $p(\mathbf{z}) = \mathcal{N}(\mathbf{0}, \mathbf{I})$
- ▶ Mapping the samples through $\mu(\mathbf{z}) = \mathbf{z}/10 + \mathbf{z}/\|\mathbf{z}\|$ yields a complex distribution
- ▶ VAEs model $\mu(\mathbf{z})$ as a **deterministic neural network** and learn its parameters

Learned Manifold



(a) Learned Frey Face manifold



(b) Learned MNIST manifold

Random Samples



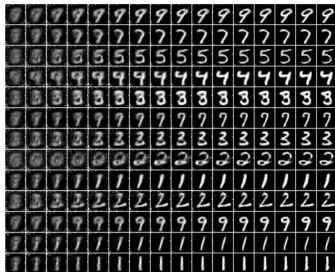
(a) 2-D latent space

(b) 5-D latent space

(c) 10-D latent space

(d) 20-D latent space

DRAW: A Sequential Variational Autoencoder



Time →

Figure 7. MNIST generation sequences for DRAW without attention. Notice how the network first generates a very blurry image that is subsequently refined.



Figure 8. Generated MNIST images with two digits.



Figure 9. Generated SVHN images.

► Sequential VAE for generating images (combines VAE with RNN)

Deep Convolutional Inverse Graphics Network

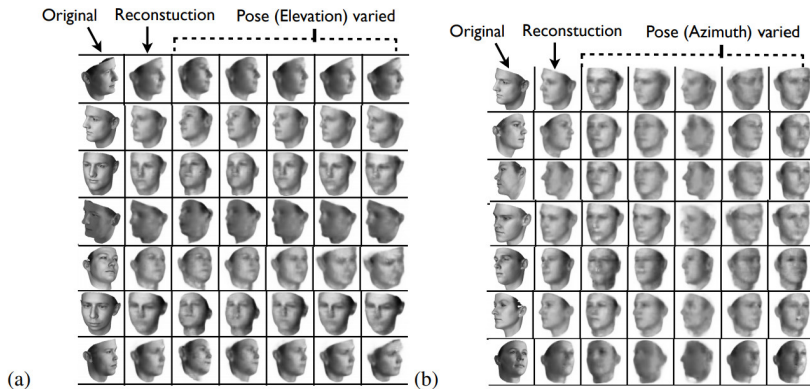
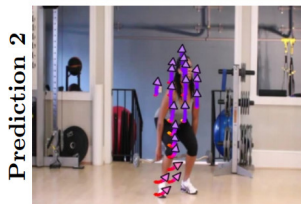


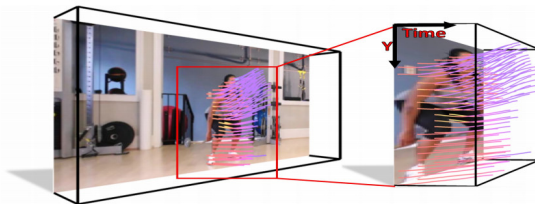
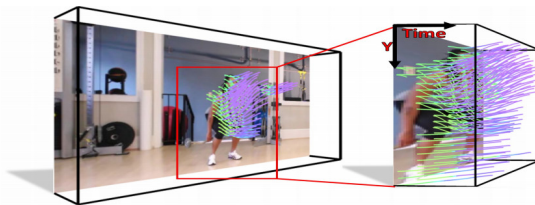
Figure 4: **Manipulating pose variables:** Qualitative results showing the generalization capability of the learned DC-IGN decoder to rerender a single input image with different pose directions.

- Forces disentangled latents (pose, light, texture, shape) through weak supervision

Motion Forecasting from Static Images



(a) Trajectories on Image



(b) Trajectories in Space-Time

- Motion forecasting from static image by jointly encoding images and trajectories

VQ-VAE-2: Vector Quantized Variational Autoencoder

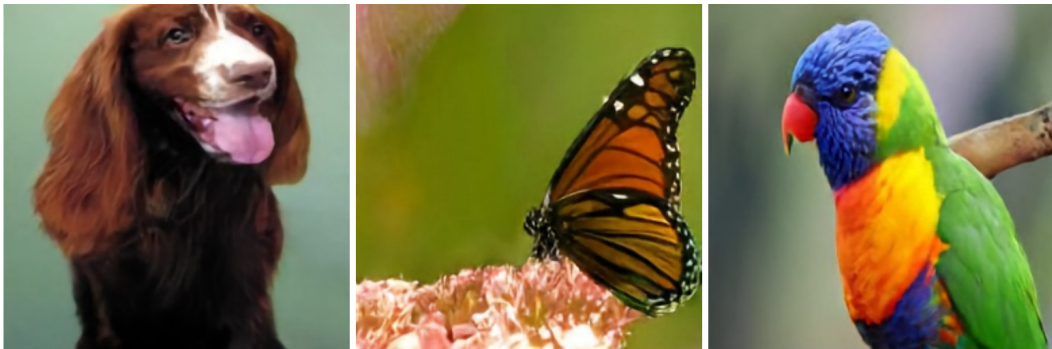
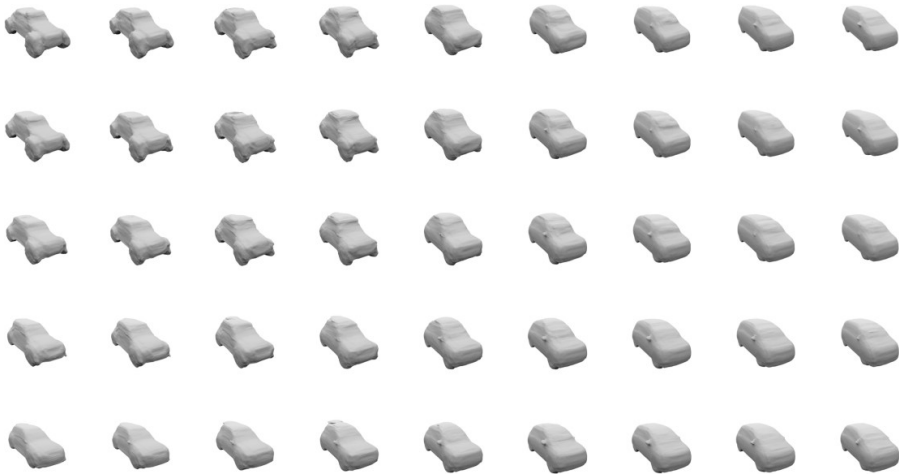


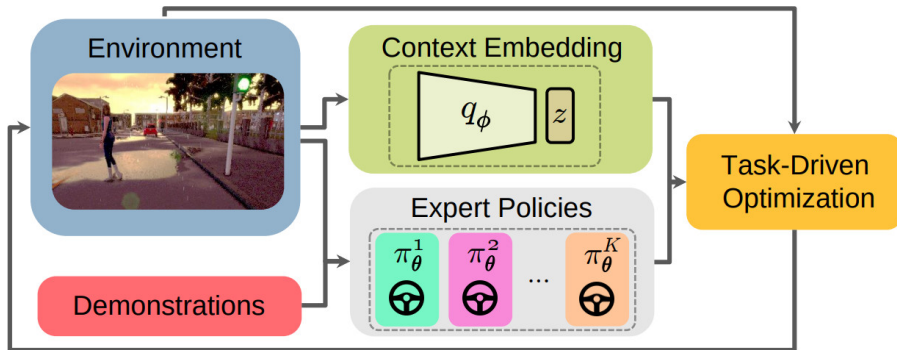
Figure 1: Class-conditional 256x256 image samples from a two-level model trained on ImageNet.

- VQ-VAEs predict discrete codes and learn the prior distribution \Rightarrow state-of-the-art

Occupancy Networks: Learning 3D Reconstruction in Function Space



Learning Situational Driving



- Data-efficient reinforcement learning with a latent embedding of the environment