Generative Models
Discriminative vs Generative

Understanding Machine Learning

Archetype: Naïve Bayes

probabilistic model:

$$P(Y|X_1,\cdots,X_p) = \frac{P(Y,X_1,\cdots,X_p)}{P(X_1,\cdots,X_p)} = \frac{P(Y)P(X_1,\cdots,X_p|Y)}{P(X_1,\cdots,X_p)} \propto P(Y)P(X_1,\cdots,X_p|Y)$$

Bayes' rule   constant   to be estimated

approach:

1. estimate $P(Y,\boldsymbol{X})$ → generative model (can be used to generate new samples)
2. calculate $P(Y|\boldsymbol{X})$ from $P(Y,\boldsymbol{X})$ → used for discriminative task (classification)

Independence Assumption

(naïve) assumption: conditional independence of features given target

$$P(X_j|Y, X_1, \cdots, X_{j-1}, X_{j+1}, \cdots, X_p) = P(X_j|Y)$$

$$\implies P(Y|X_1, \cdots, X_p) = \frac{P(Y)\prod_{j=1}^{p} P(X_j|Y)}{P(X_1, \cdots, X_p)}$$
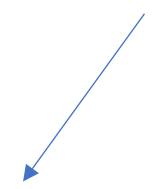
→ independent feature contributions (ignoring feature correlations)
→ robust against curse of dimensionality

Estimation of Feature Contributions

separate estimations of $P(X_j|Y)$ for each feature

requires assumption of distributions (e.g., Gaussian naïve Bayes) or non-parametric methods (kernel density estimation)

Gaussian feature likelihoods: $P(x_{ij}|y) = \dfrac{1}{\sqrt{2\pi\sigma_{y,j}^2}} \exp\left(-\dfrac{(x_{ij}-\mu_{y,j})^2}{2\sigma_{y,j}^2}\right)$

parameter estimation (e.g., mean and variance of Gaussians) can be done with maximum likelihood method (y known in training)

$\rightarrow$ no Bayesian methods needed

Maximum a Posteriori Classification

$$\hat{y}_i = \underset{y}{\text{argmax}} \, P(y) \prod_{j=1}^{p} P(x_{ij}|y)$$

despite potentially inaccurate probability estimates (due to naïve independence assumption), good identification of correct class via maximum probability

→ bad for regression tasks (if independence assumption is too naïve, i.e., features are correlated)
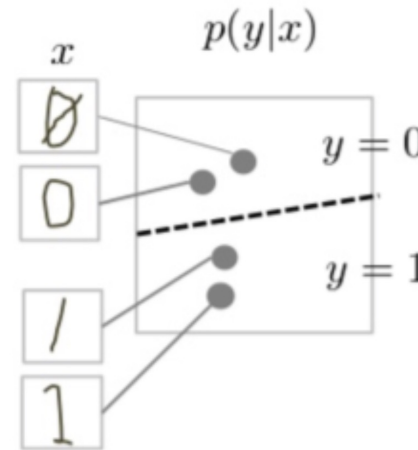
Generative vs Discriminative Models

generative models: predict joint probability $P(Y, \boldsymbol{X})$ (what allows to create new data samples) or directly generates new data samples
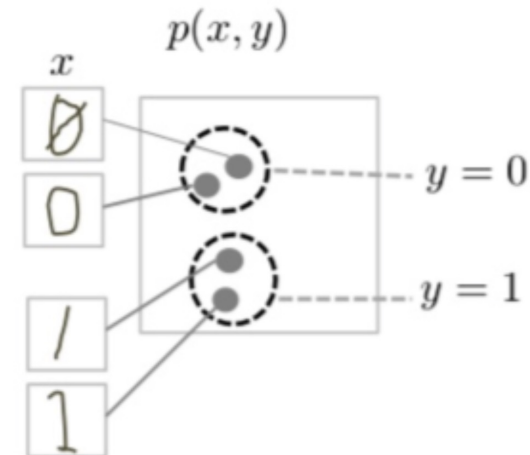
discriminative models: predict conditional probability $P(Y|\boldsymbol{X})$ or directly output (label for classification, real value for regression)

task of generative models more difficult: model full data distribution rather than merely find patterns in inputs to distinguish outputs

discriminative model

generative model

Naïve Bayes and Logistic Regression

generative-discriminative pair of classification algorithms

- binary case: logit of naïve Bayes' outputs, $\log\left(\frac{P(y_i=1|\boldsymbol{x}_i)}{P(y_i=0|\boldsymbol{x}_i)}\right)$, corresponds to output of logistic regression's linear predictor
- for discrete inputs or Gaussian naïve Bayes: naïve Bayes can be reparametrized as linear classifier

for discriminative task: identical in asymptotic limit (infinite training samples) if independence assumption holds (otherwise naïve Bayes less accurate)

naïve Bayes has greater bias but lower variance than logistic regression $\rightarrow$ to be preferred for scarce training data (if bias, i.e., independence assumption, correct)
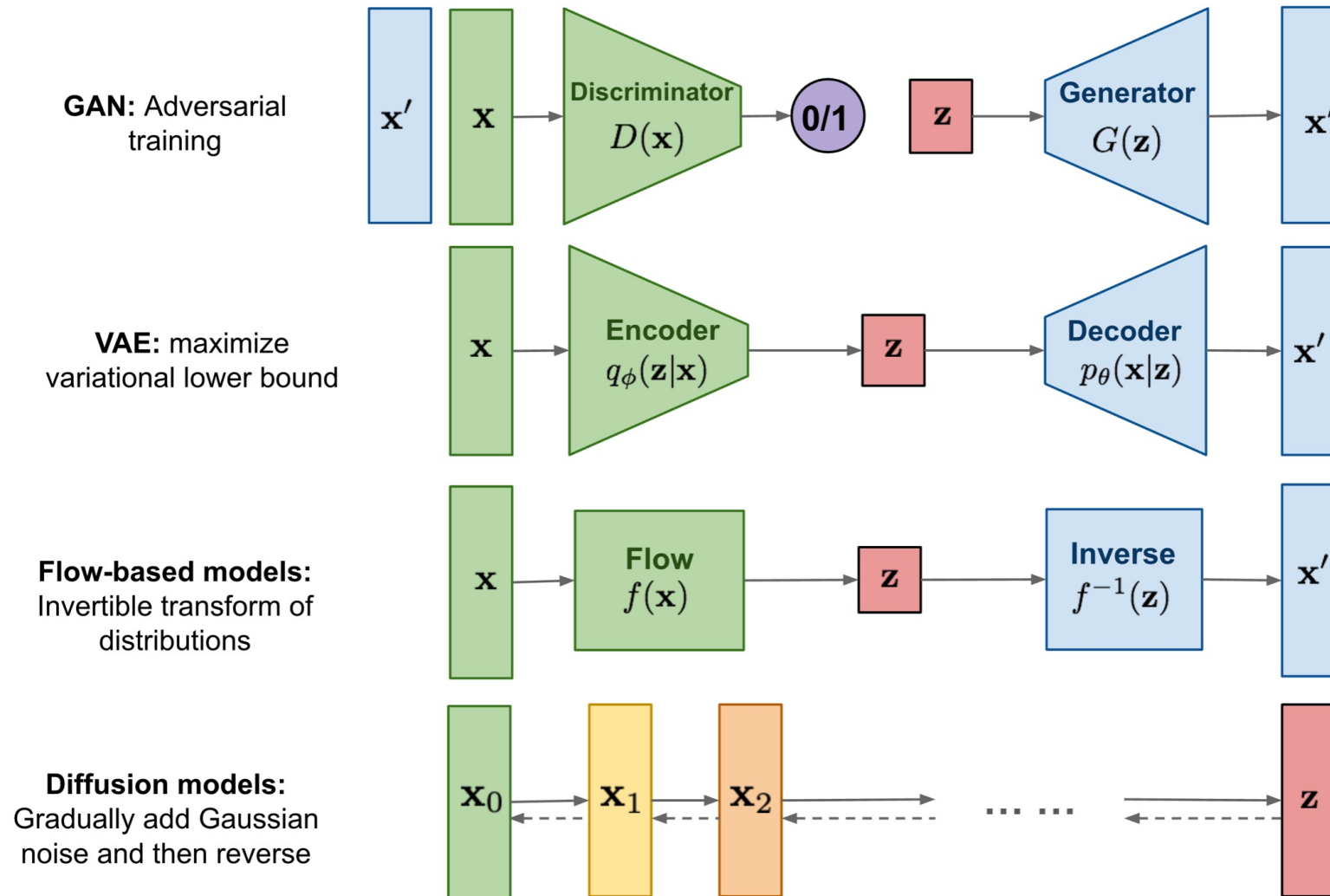
Data Generation

generative models can be used for discriminative tasks (although potentially inferior to direct discriminative methods)

but generative methods do more than discriminative ones: model full data distribution

→ allows generation of new data samples (can be images, text, video, audio, proteins, materials, time series, structured data, …)

large (auto-regressive) language models examples of generative models

Different Types of Generative Models

GAN: Adversarial training

VAE: maximize variational lower bound

Flow-based models: Invertible transform of distributions

Diffusion models: Gradually add Gaussian noise and then reverse
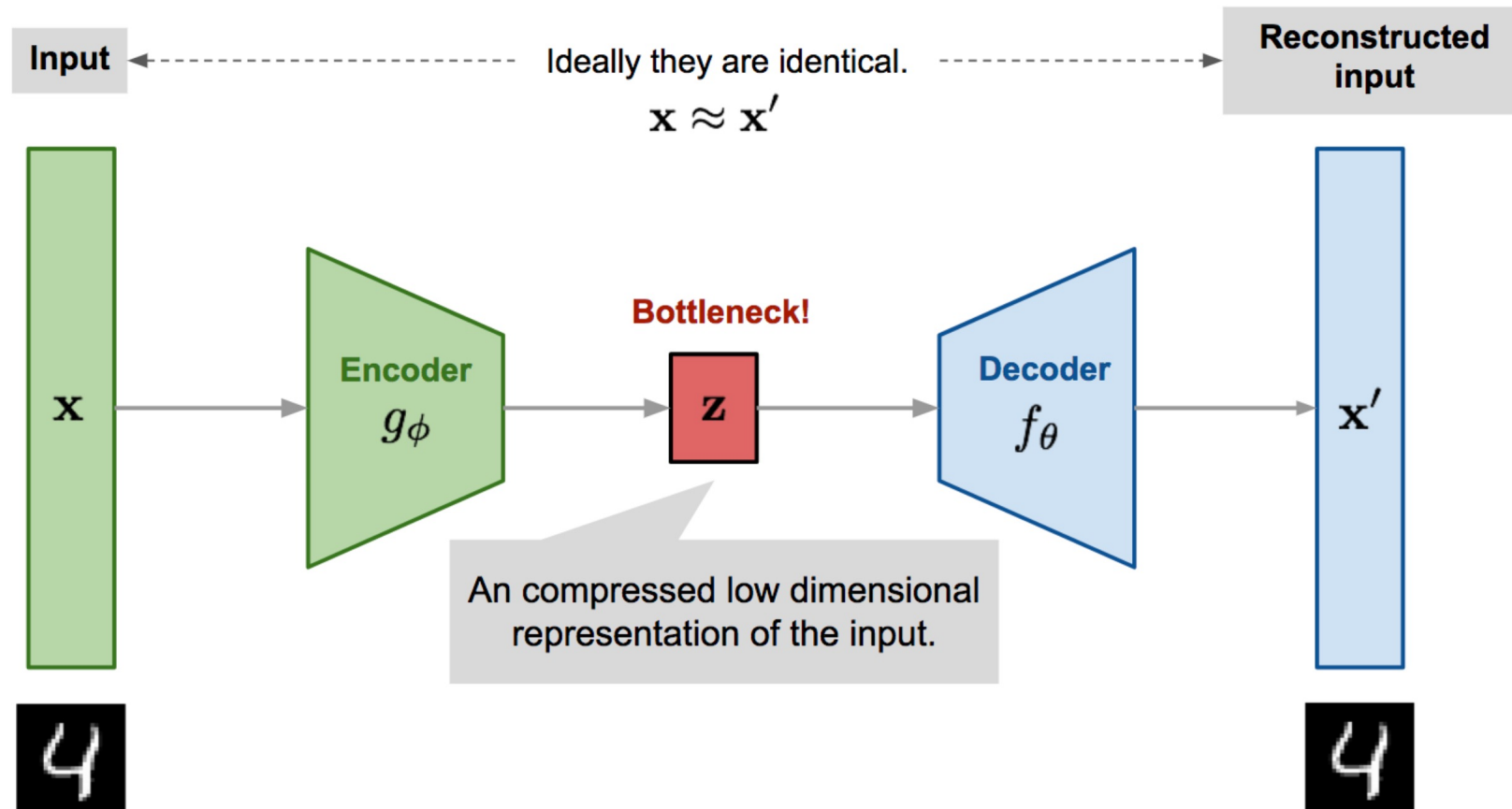
source

Variational Autoencoders (VAE)

Recap: Autoencoder

(deep) encoder network

(deep) decoder network

learned together by minimizing differences between original input and reconstructed input (expressed as losses)

compressed intermediate representation: dimensionality reduction

Input — — — — — — — — Ideally they are identical. — — — — — — Reconstructed input

$$\mathbf{x} \approx \mathbf{x}'$$

Bottleneck!

Encoder g_ϕ

$\mathbf{z}$

Decoder f_θ

$\mathbf{x}$

$\mathbf{x}'$

An compressed low dimensional representation of the input.

Autoencoder Architecture for Generative Tasks

goal: generation of variations of input data rather than compressed representation

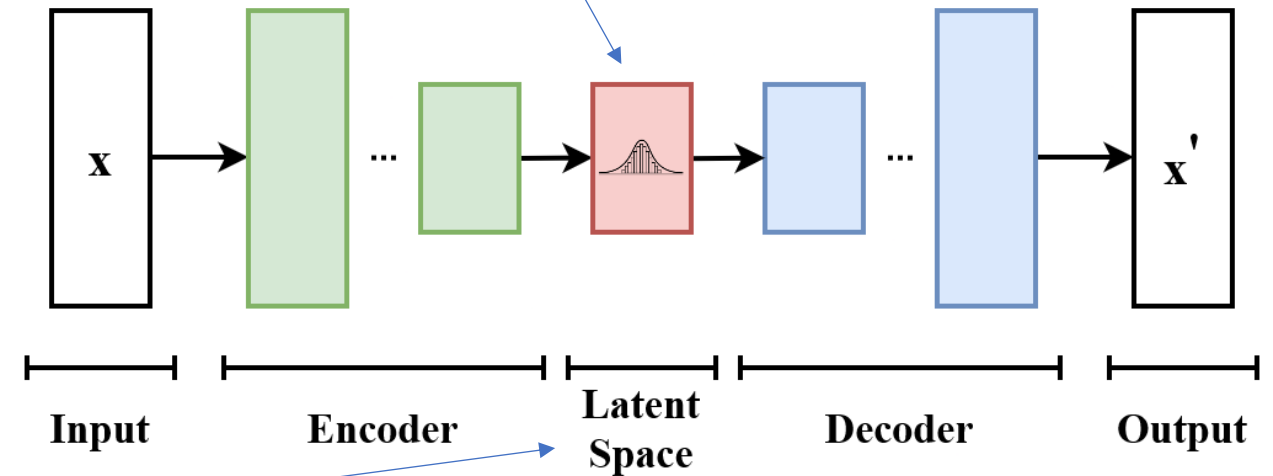→ learn variational distribution instead of identity function

to be precise: parametrized variational distribution of latent encoding variables $\boldsymbol{z}$

prior (simple distribution, in usual VAE: Gaussian): $p_{\boldsymbol{\theta}}(\boldsymbol{z})$

posterior: $p_{\boldsymbol{\theta}}(\boldsymbol{z}|\boldsymbol{x}) = \dfrac{p_{\boldsymbol{\theta}}(\boldsymbol{x}|\boldsymbol{z})p_{\boldsymbol{\theta}}(\boldsymbol{z})}{\int p_{\boldsymbol{\theta}}(\boldsymbol{x}|\boldsymbol{z})p_{\boldsymbol{\theta}}(\boldsymbol{z})d\boldsymbol{z}}$

$p_{\boldsymbol{\theta}}(\boldsymbol{x})$: mixture of Gaussians

from which to sample

| Input | Encoder | Latent Space | Decoder | Output |

from wikipedia

Variational Bayesian Method

12

Encoder and Decoder Networks

encoder: find posterior $p_{\boldsymbol{\theta}}(\boldsymbol{z}|\boldsymbol{x})$

unfortunately, generally intractable (integral over $\boldsymbol{z}$ expensive)

$\rightarrow$ approximate by $q_{\boldsymbol{\phi}}(\boldsymbol{z}|\boldsymbol{x})$

VAE: $q_{\boldsymbol{\phi}}(\boldsymbol{z}|\boldsymbol{x})$ expressed by neural network with weights $\boldsymbol{\phi}$

$\rightarrow$ amortized inference:
$q_{\boldsymbol{\phi}}(\boldsymbol{z}|\boldsymbol{x})$ learned in training, $\boldsymbol{z}$ inferred from $\boldsymbol{x}$ in prediction (sharing variational parameters across all data points)

in VAE: network weights $\boldsymbol{\theta}$

decoder: generate new sample $\boldsymbol{x}_i$

1. sample $\boldsymbol{z}_i$ (from Gaussian)

2. generate $\boldsymbol{x}_i$ (similar to real data)

$\rightarrow$ maximize: $p_{\boldsymbol{\theta}}(\boldsymbol{x}_i) = \int p_{\boldsymbol{\theta}}(\boldsymbol{x}_i|\boldsymbol{z})\, p_{\boldsymbol{\theta}}(\boldsymbol{z}) d\boldsymbol{z}$

(expensive $\rightarrow$ use only likely codes $\boldsymbol{z}$ given input $\boldsymbol{x}$: need for encoder)

encoder network

decoder network

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

$p_\theta(\mathbf{z})$
$\mathbf{z} \sim \mathcal{N}(0,1)$

$\mathbf{z}$

$p_\theta(\mathbf{x}|\mathbf{z})$

$\mathbf{x}$

$p_\theta(\mathbf{x})$

13
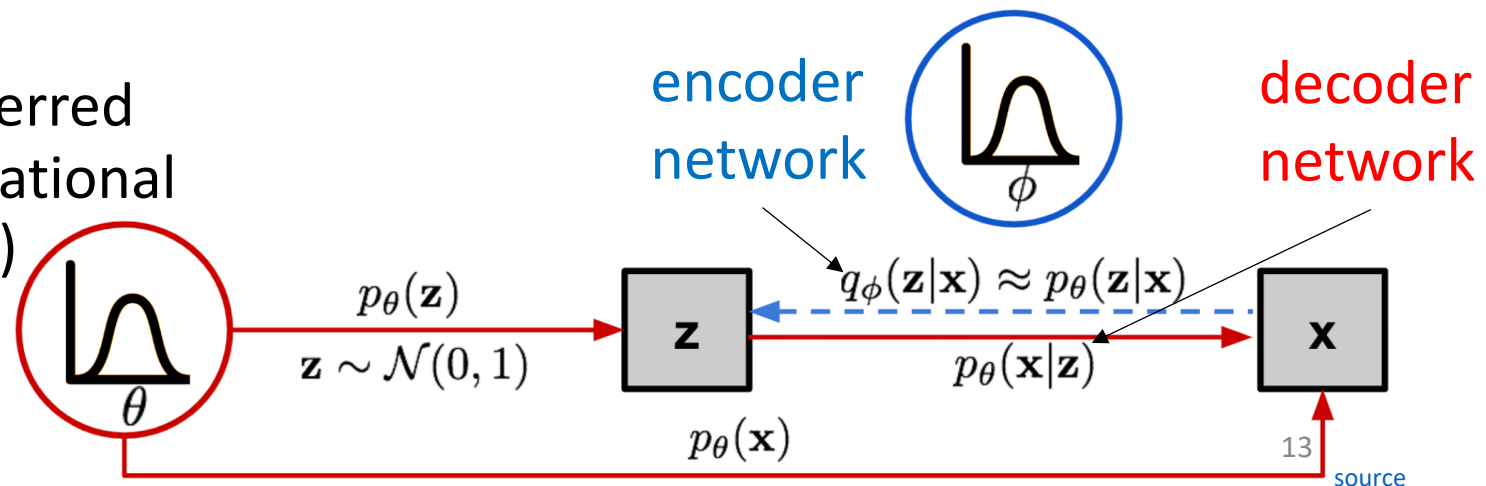
VAE Loss: ELBO

VAE loss function to be minimized according to network weights:

$$L(\boldsymbol{x}_i; \boldsymbol{\theta}, \boldsymbol{\phi}) = -\ln p_{\boldsymbol{\theta}}(\boldsymbol{x}_i) + D_{KL}\left(q_{\boldsymbol{\phi}}(\boldsymbol{z}|\boldsymbol{x}_i)||p_{\boldsymbol{\theta}}(\boldsymbol{z}|\boldsymbol{x}_i)\right)$$

maximize likelihood of observed data (minimize reconstruction error)

and

minimize difference of approximation $q_{\boldsymbol{\phi}}(\boldsymbol{z}|\boldsymbol{x}_i)$ to exact posterior $p_{\boldsymbol{\theta}}(\boldsymbol{z}|\boldsymbol{x}_i)$

can be interpreted as regularizer

corresponds to maximizing evidence lower bound (ELBO), i.e., maximizing lower bound of probability to generate real data sample:
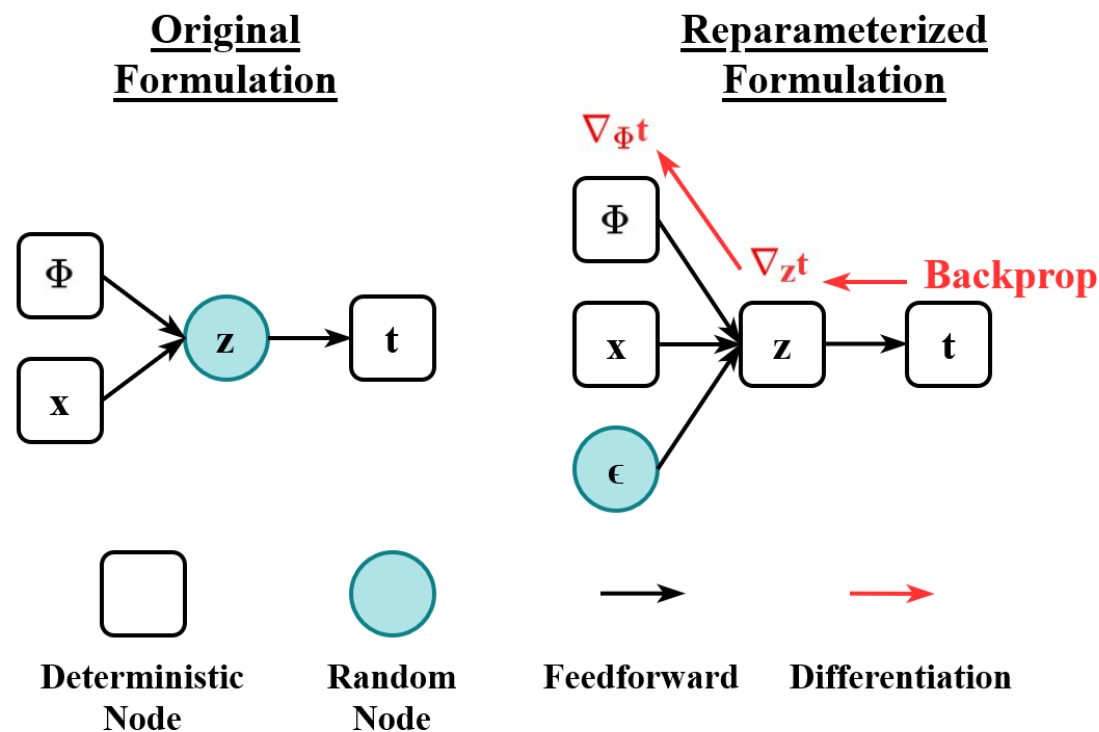
$$\ln p_{\boldsymbol{\theta}}(\boldsymbol{x}_i) \geq \ln p_{\boldsymbol{\theta}}(\boldsymbol{x}_i) - \underbrace{D_{KL}\left(q_{\boldsymbol{\phi}}(\boldsymbol{z}|\boldsymbol{x}_i)||p_{\boldsymbol{\theta}}(\boldsymbol{z}|\boldsymbol{x}_i)\right)}_{\text{non-negative}} = E_{\boldsymbol{z}\sim q_{\boldsymbol{\phi}}(\boldsymbol{z}|\boldsymbol{x}_i)}\left[\ln \frac{p_{\boldsymbol{\theta}}(\boldsymbol{x}_i, \boldsymbol{z})}{q_{\boldsymbol{\phi}}(\boldsymbol{z}|\boldsymbol{x}_i)}\right]$$

Reparameterization Trick

→ gradient descent according to $\boldsymbol{\theta}$ and $\boldsymbol{\phi}$

issue: not readily possible for $\boldsymbol{\phi}$
(expecatation over $\boldsymbol{z}$, which is sampled
from $q_{\boldsymbol{\phi}}$)

→ reparametrization to the rescue:
express randomness in $\boldsymbol{z}$ by independent
auxiliary variable $\boldsymbol{\varepsilon}$

from wikipedia

e.g., q_ϕ as multivariate Gaussian with diagonal covariance structure
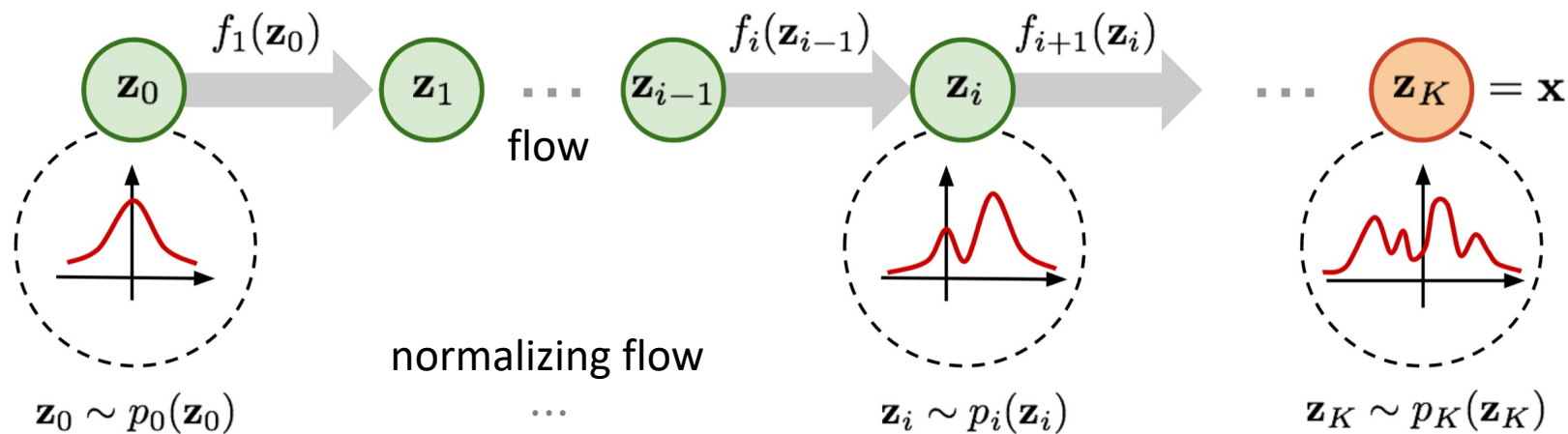
$\rightarrow$ learn mean and variance
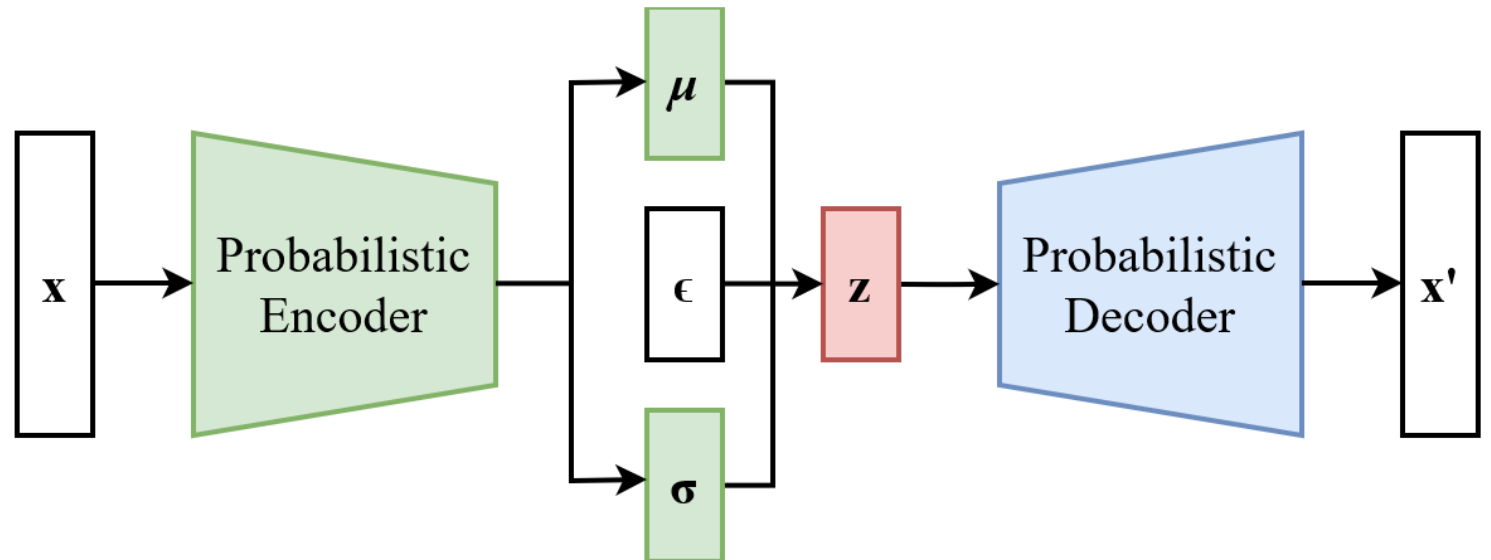
Flow-Based Methods

Normalizing Flows

idea: mapping of a simple probability distribution (often, standard normal distribution) into a complex one by sequence of invertible transformations (repeatedly applying the change-of-variable technique)

$$f(\mathbf{z}') = f(\mathbf{z}) \left| \det \frac{\delta f^{-1}}{\delta \mathbf{z}'} \right| = f(\mathbf{z}) \left| \det \frac{\delta f}{\delta \mathbf{z}} \right|^{-1}$$

log-likelihood:
$$\ln p_K(\mathbf{z}_K) = \ln p_0(\mathbf{z}_0) - \sum_{k=1}^{K} \ln \left| \det \frac{\delta f_k}{\delta \mathbf{z}_{k-1}} \right|$$

Usage in Generative Models

training: estimate maximum likelihood of normalizing flow (log-likelihood of last slide) by gradient descent (learn parameters θ of transformations f_θ^{-1}, e.g., to let $p_0(\boldsymbol{z})$ be Gaussian)

inference: sample from simple distribution $p_0(\boldsymbol{z})$ and transform it back to data distribution $p_K(\boldsymbol{x})$ via f_θ

advantages over VAE:

- instead of simple functions like Gaussians, allow more complex ones: real-world distributions usually much more complicated

- direct/explicit estimation of likelihood (negative log-likelihood as loss): allows density estimation (e.g., to predict rareness of future events)

Invertible Neural Networks

neural networks representing invertible/bijective functions can be used for normalizing flow transformations

need for specialized architectures to construct reversible transform (e.g., affine coupling layers)

Generative Adversarial Networks (GAN)

Indirect Training via Discriminator

two neural networks playing a zero-sum game:

- the generator network G generating new (fake) samples
- the discriminator network D trying to distinguish between real and fake samples

idea: G not trained directly to minimize reconstruction error of real samples, but to fool D → self-supervised approach

Formulation
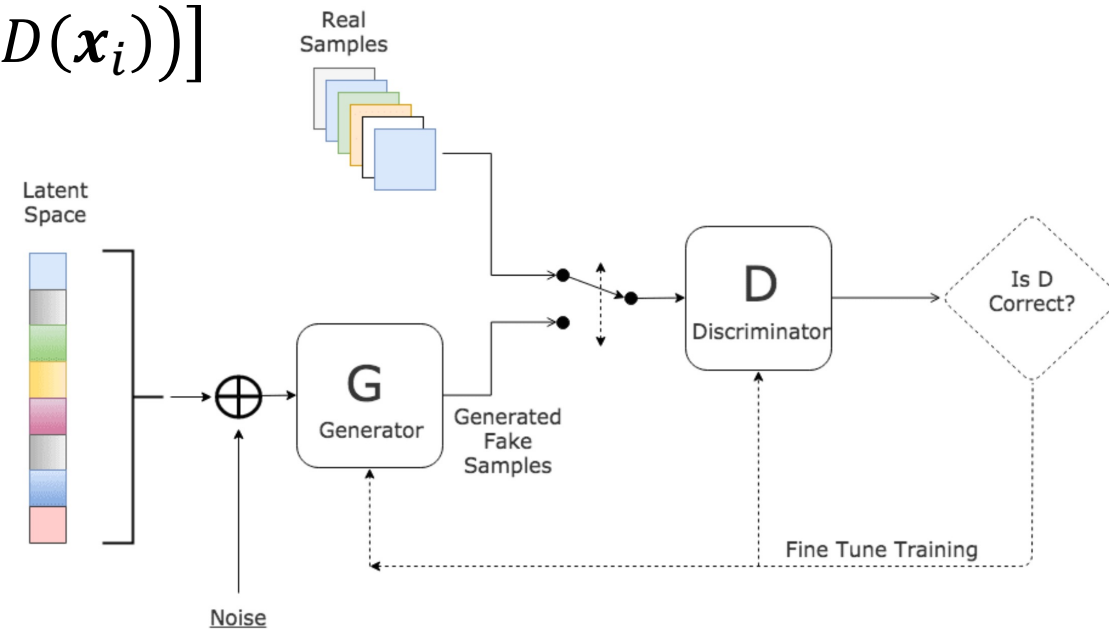
common loss for generator and discriminator:

$$L(\boldsymbol{x_i}) = E_{\boldsymbol{x} \sim p_r(\boldsymbol{x})}[\ln D(\boldsymbol{x_i})] + E_{\boldsymbol{x} \sim p_g(\boldsymbol{x})}\left[\ln\big(1 - D(\boldsymbol{x_i})\big)\right]$$

- G trying to minimize
- D trying to maximize

decomposition into latent space (parameters of generator network) and noise (sampled from, e.g., Gaussian distribution): reparametrization trick

source

23

Properties

implicit generative model: do not estimate likelihood function

for optimal D, GAN loss quantifies similarity between generative data distribution p_g and real data distribution p_r by Jensen-Shannon divergence

$$D_{JS}(p||q) = \frac{1}{2}D_{KL}\left(p||\frac{p+q}{2}\right) + \frac{1}{2}D_{KL}\left(q||\frac{p+q}{2}\right)$$

for optimal values of both G and D: $p_g = p_r$ and $D = 0.5$

issue: potentially unstable training

Diffusion Models

Idea

training:

distort training data by successively adding random noise, then learn to reverse this process (denoising)

generation:

sample random noise and run through the learned denoising process

can be seen as special kind of energy-based models: parametrized energy function (to be minimized) of target, observable, and latent variables → no need for proper normalization of probability distributions

Forward Process

Markov chain of diffusion steps to slowly add Gaussian noise to data (inspired by non-equilibrium thermodynamics):

$$q(\boldsymbol{x}_{1:T}|\boldsymbol{x}_0) = \prod_{t=1}^{T} q(\boldsymbol{x}_t|\boldsymbol{x}_{t-1}), \quad q(\boldsymbol{x}_t|\boldsymbol{x}_{t-1}) = \mathcal{N}\left(\boldsymbol{x}_t; \sqrt{1-\beta_t}\boldsymbol{x}_{t-1}, \beta_t \mathbf{I}\right)$$

- with variance schedule $\beta_1, \ldots, \beta_T$ (hyperparameters, increasing with t)
- large T and small β_t → same functional form for forward and reverse processes, ending up with isotropic Gaussian distribution for $\boldsymbol{x}_T$

$$t = 0 \qquad\qquad t = \frac{T}{2} \qquad\qquad t = T$$

$q\left(\mathbf{x}^{(0\cdots T)}\right)$
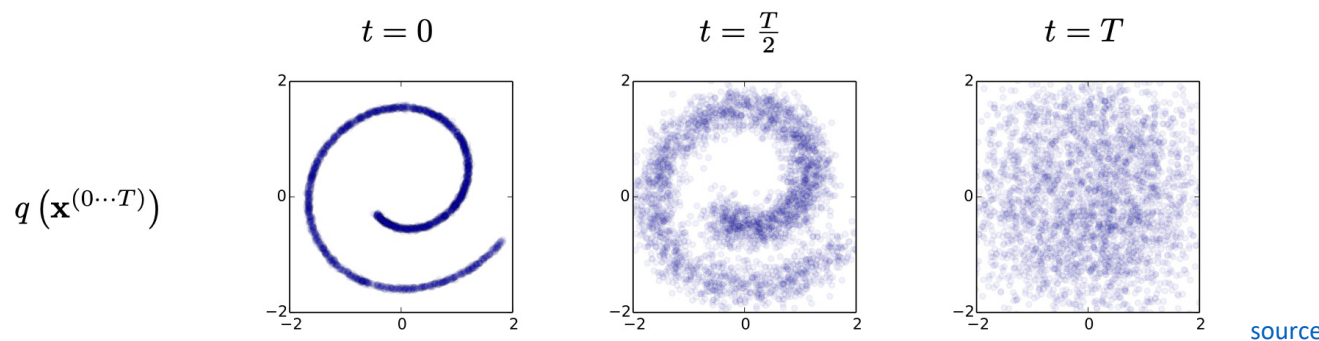
Reparametrization

conditional Gaussian distributions at each t:

sample $\epsilon \sim \mathcal{N}(0, \mathbf{I})$ and set $\boldsymbol{x}_t = \sqrt{1 - \beta_t} \boldsymbol{x}_{t-1} + \sqrt{\beta_t} \epsilon$

nice property: possible to directly sample $\boldsymbol{x}_t$ conditioned on $\boldsymbol{x}_0$ (no need to apply q repeatedly)

$$\boldsymbol{x}_t = \sqrt{\bar{\alpha}_t} \boldsymbol{x}_0 + \sqrt{1 - \bar{\alpha}_t} \epsilon$$
$$q(\boldsymbol{x}_t | \boldsymbol{x}_0) = \mathcal{N}\left(\boldsymbol{x}_t; \ \sqrt{\bar{\alpha}_t} \boldsymbol{x}_0, (1 - \bar{\alpha}_t)\mathbf{I}\right)$$

with $\quad \alpha_t = 1 - \beta_t, \quad \bar{\alpha}_t = \prod_{s=1}^{t} \alpha_s$

conditioning on $\boldsymbol{x}_0$ also allows to handle $q(\boldsymbol{x}_{t-1} | \boldsymbol{x}_t, \boldsymbol{x}_0)$

Reverse Process

to generate new data samples, one needs to learn to reverse the diffusion process (starting from pure noise): neural network learning to gradually denoise data

overall loss as sum of losses for each time step t

for each t: D_{KL} between two Gaussians (closed form) $q(\boldsymbol{x}_{t-1}|\boldsymbol{x}_t, \boldsymbol{x}_0)$ and $p_\theta(\boldsymbol{x}_{t-1}|\boldsymbol{x}_t)$
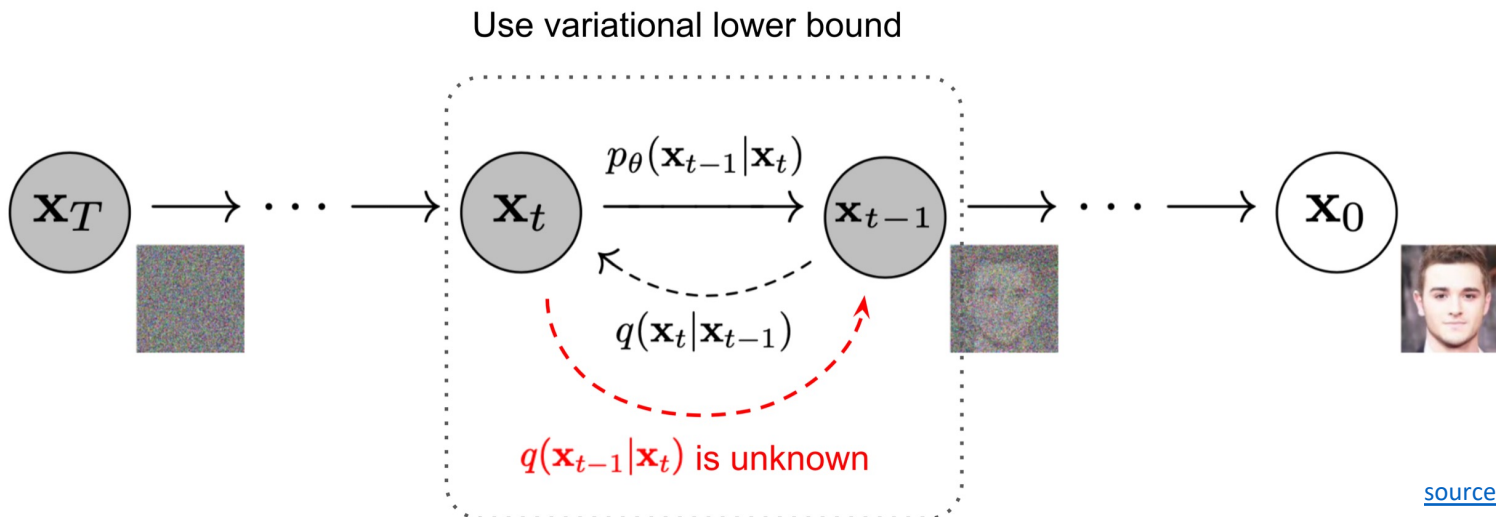
→ corresponds to VAE loss: maximizing ELBO

time-dependent Gaussian parameters:
$$p(\boldsymbol{x}_T) = \mathcal{N}(\boldsymbol{x}_T; 0, \mathbf{I})$$

$$p_\theta(\boldsymbol{x}_{0:T}) = p(\boldsymbol{x}_T) \prod_{t=1}^{T} p_\theta(\boldsymbol{x}_{t-1}|\boldsymbol{x}_t)$$

$$p_\theta(\boldsymbol{x}_{t-1}|\boldsymbol{x}_t) = \mathcal{N}(\boldsymbol{x}_{t-1}; \boldsymbol{\mu}_\theta(\boldsymbol{x}_t, t), \boldsymbol{\Sigma}_\theta(\boldsymbol{x}_t, t))$$

Use variational lower bound

$p_\theta(\mathbf{x}_{t-1}|\mathbf{x}_t)$

$\mathbf{x}_T \longrightarrow \cdots \longrightarrow \mathbf{x}_t \longrightarrow \mathbf{x}_{t-1} \longrightarrow \cdots \longrightarrow \mathbf{x}_0$

$q(\mathbf{x}_t|\mathbf{x}_{t-1})$

$q(\mathbf{x}_{t-1}|\mathbf{x}_t)$ is unknown

Noise Prediction

reparametrization allows to learn added noise instead of Gaussian parameters:

Algorithm 1 Training

1: **repeat**
2:   $\mathbf{x}_0 \sim q(\mathbf{x}_0)$
3:   $t \sim \text{Uniform}(\{1, \ldots, T\})$
4:   $\boldsymbol{\epsilon} \sim \mathcal{N}(\mathbf{0}, \mathbf{I})$
5:   Take gradient descent step on
$$\nabla_\theta \left\| \boldsymbol{\epsilon} - \boldsymbol{\epsilon}_\theta(\sqrt{\bar{\alpha}_t}\mathbf{x}_0 + \sqrt{1 - \bar{\alpha}_t}\boldsymbol{\epsilon}, t) \right\|^2$$
6: **until** converged

Algorithm 2 Sampling

1: $\mathbf{x}_T \sim \mathcal{N}(\mathbf{0}, \mathbf{I})$
2: **for** $t = T, \ldots, 1$ **do**
3:   $\mathbf{z} \sim \mathcal{N}(\mathbf{0}, \mathbf{I})$ if $t > 1$, else $\mathbf{z} = \mathbf{0}$
4:   $\mathbf{x}_{t-1} = \frac{1}{\sqrt{\alpha_t}}\left(\mathbf{x}_t - \frac{1 - \alpha_t}{\sqrt{1 - \bar{\alpha}_t}}\boldsymbol{\epsilon}_\theta(\mathbf{x}_t, t)\right) + \sigma_t \mathbf{z}$
5: **end for**
6: **return** $\mathbf{x}_0$

source

L2-loss (MSE) between true and predicted Gaussian noise at time step t

use position embeddings (as network parameters are shared across time)

diffusion models can be interpreted as chain of denoising autoencoders (also connected to score-based generative modeling via Langevin dynamics)

Denoising Autoencoder

goal: avoid overfitting and improve robustness of plain autoencoder

learn to remove noise of distorted input $\tilde{x}$ → restore original input x

similar to dropout
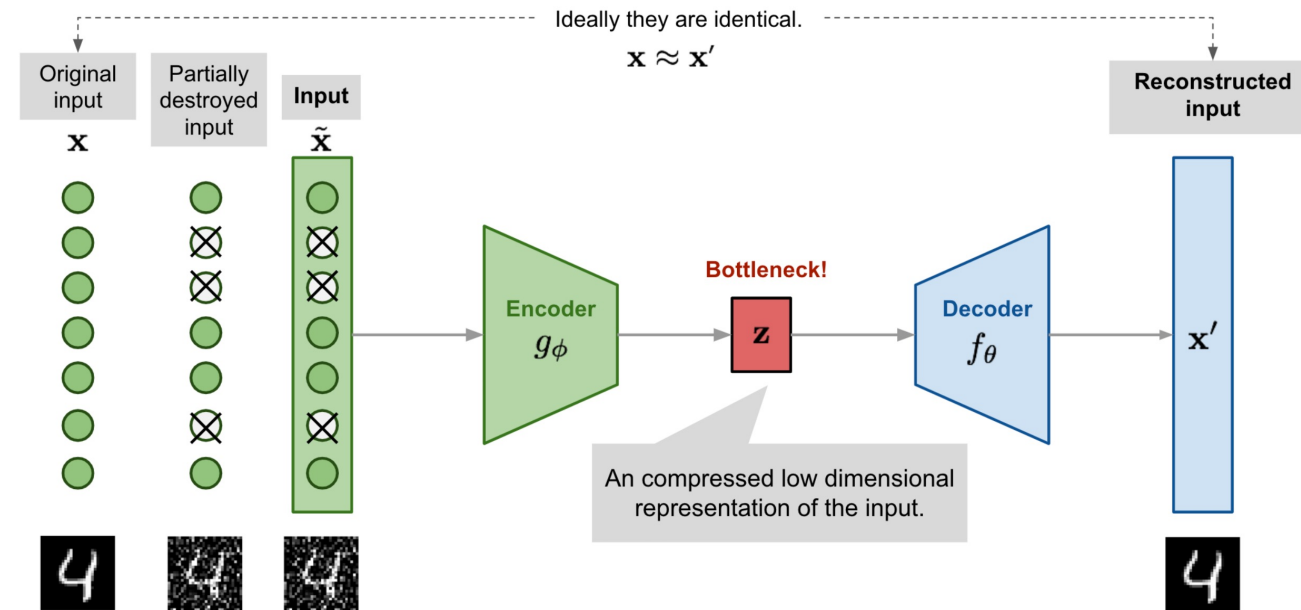
differences of diffusion models to typical denoising autoencoders:
- no bottleneck (care about output here, not internal representation): latent space with high dimensionality (same as original data)
- handle many different noise levels with single set of shared parameters

Latent Diffusion Model

add noise to latent representation rather than raw data

→ significant speedup

diffusion model highly flexible in terms of architecture: only require same input and output dimensionality → often U-Net-like architectures (autoencoders) with skip connections

source

32
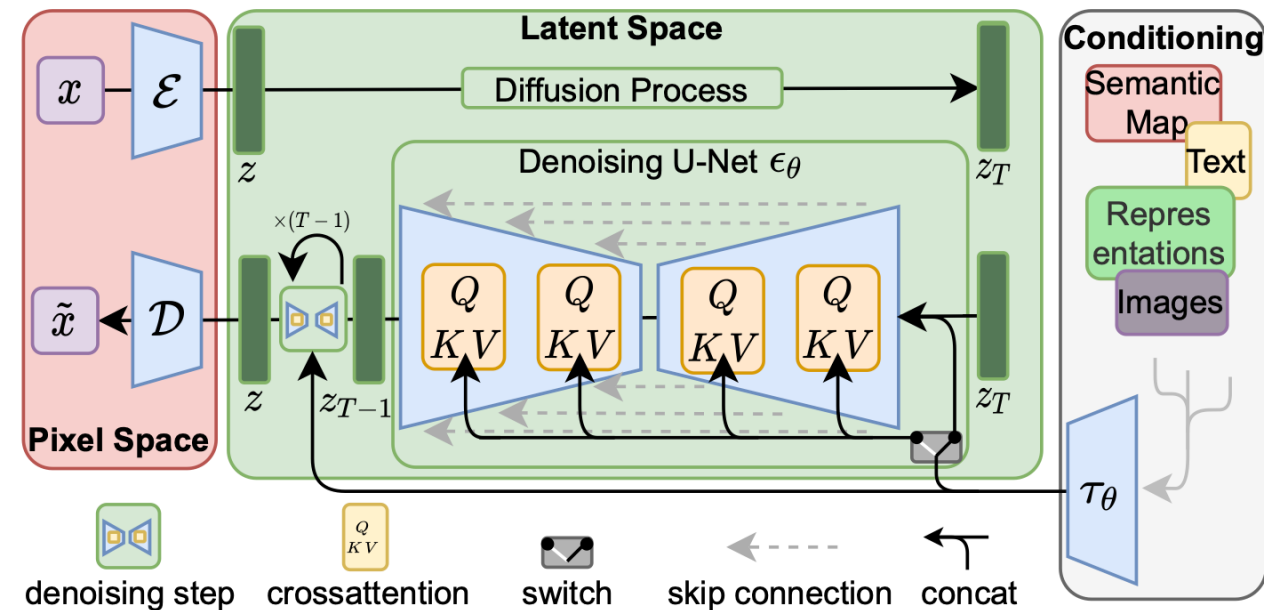
Conditioned Generation

as discussed so far, generative methods give no control over what kind of data is generated (limited usability)

→ need for conditional approach (e.g., conditioning on describing text)

… attention

classifier-guided diffusion: …

classifier-free guidance: …

Applications

plenty of implementations available

text-to-image:
[DALL-E 2](), [Stable Diffusion](), [GLIDE](),
[ImageGen]()

web app for Stable Diffusion:

[DreamStudio]()

A dream of a distant galaxy, by Caspar David Friedrich, matte painting trending on artstation HQ

Literature

papers:

- variational autoencoder

- normalizing flows

- GAN

- denoising diffusion, latent diffusion

Movie-like Intelligence

emergent capabilities of complex systems almost impossible to foresee

mini examples in contemporary ML:

- large language models
- multi-agent reinforcement learning

one idea: reward is enough

philosophical: emotions or consciousness might also occur as emergent capabilities