CS/DS 541: Class 19

Jacob Whitehill

Exercises

https://cs230.stanford.edu/files/cs230exam_fall19_soln.pdf

Exercise 4

Consider a model trying to learn an encoding of some input $x \in \mathbb{R}$. The goal is to encode the input x using $z = w_1 x \in \mathbb{R}$, then accurately reconstruct the original x from the encoded representation using $\hat{x} = w_2 z \in \mathbb{R}$. Here, $(w_1, w_2) \in \mathbb{R} \times \mathbb{R}$. The model is trained with the squared reconstruction error:

$$L(W) = \frac{1}{n} \sum_{i=1}^{n} (x^{(i)} - w_2 w_1 x^{(i)})^2$$

(2 points) What is the set of solutions for w_1 and w_2 which makes loss zero?

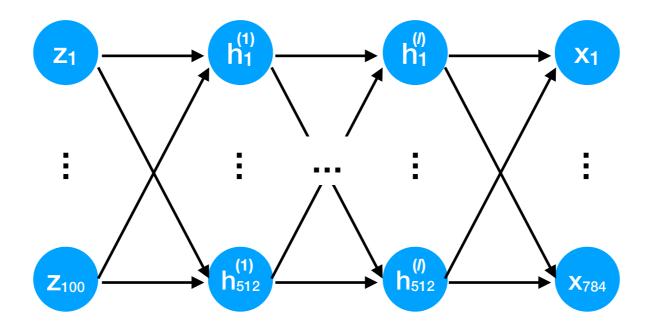
(3 points) Does the loss have a saddle point? Where?

- However, another entire class of deep learning methods is based on training two networks that compete against each other in a zero-sum game.
- In particular, the most prominent method (as of 2020) for generating novel data x is the Generative Adversarial Network (GAN; Goodfellow et al. 2014).

- Like VAEs, GANs consist of two components, but their semantics are different.
- Let $P_{\text{data}}(\mathbf{x})$ be the ground-truth data distribution.
- Generator G: given a noise vector z from an easy-to-sample distribution (e.g., Gaussian, uniform), generate a vector x that looks like it came from P_{data}(x).
- **Discriminator** D: given a vector x, decide if it is real ($\hat{y}=1$) or fake ($\hat{y}=0$). D acts as a "forgery detector".

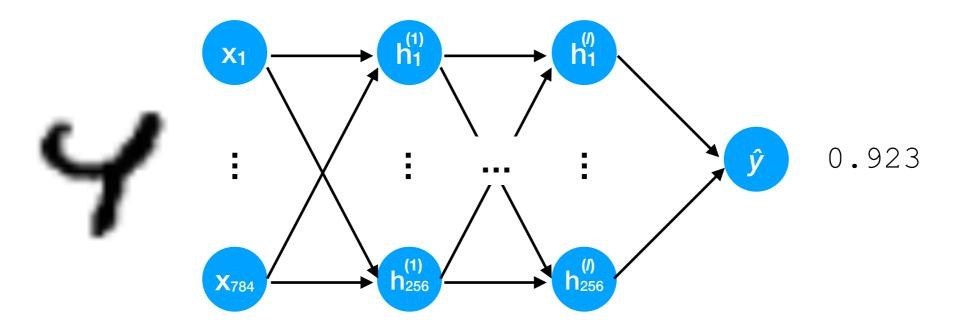
Generator G

Example G with I hidden layers that generates an MNIST image (28x28=784) x from a 100-dim noise vector z:

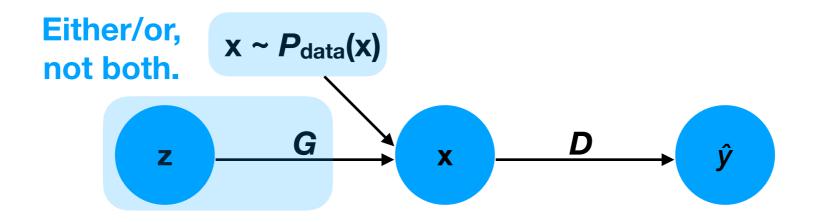


Discriminator D

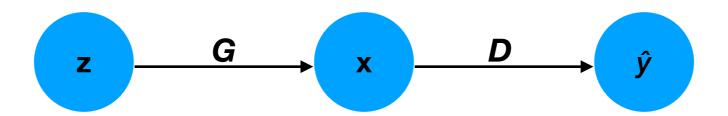
• Example D with I hidden layers that estimates $\hat{y} \in (0,1)$ that expresses probability that the input \mathbf{x} is real:



- Like VAEs, GANs are trained such that one component "feeds" to the other.
- In contrast to VAEs, the discriminator D is sometimes given a "fake" data vector \mathbf{x} (generated by G), and sometimes given a "real" vector \mathbf{x} sampled from the training set (which approximates $P_{\text{data}}(\mathbf{x})$).



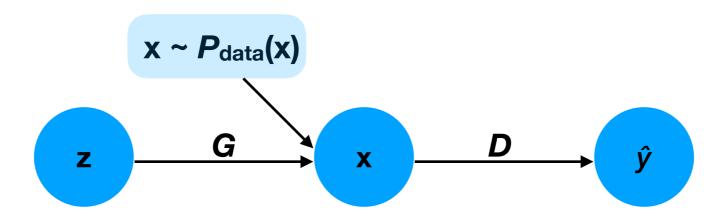
- Each network has its own parameters:
 - G has parameters θ_G .
 - *D* has parameters θ_D .



 We can define the following loss on how well D can discriminate fake from real data:

$$f_{\text{acc}}(\theta_G, \theta_D) = \mathbb{E}_{\mathbf{x} \sim P_{\text{data}}(\mathbf{x})} [\log D_{\theta_D}(\mathbf{x})] + \mathbb{E}_{\mathbf{z} \sim P(\mathbf{z})} [\log (1 - D_{\theta_D}(G_{\theta_G}(\mathbf{z})))]$$

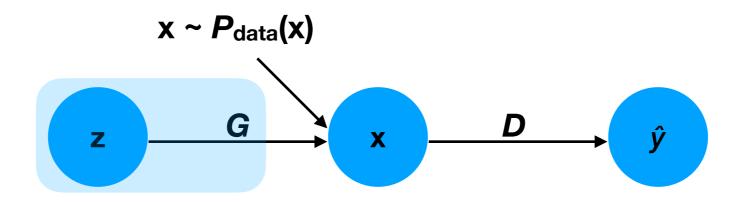
Log-likelihood that *D* recognizes real data as real.



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Log-likelihood that *D* recognizes fake data as fake.



• The goal of D is to maximize f_{acc} , whereas the goal of G is to minimize f_{acc} .

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This two-player game will reach an equilibrium if we find:

$$\min_{\theta_G} \max_{\theta_D} f_{\rm acc}(\theta_G, \theta_D)$$

In particular, this solution corresponds to D having 50% accuracy at detecting forgeries, and G generating fake x according to P_{data}(x).

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- In practice, we train *D* and *G iteratively*:
 - Freeze G, and perform SGD on D for k iterations to increase f_{acc}.

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Improve *D*'s forgery detection accuracy for a fixed distribution of fake data.

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 - Freeze G, and perform SGD on D for k iterations to increase f_{acc}.
 - Freeze D, and perform SGD on G for I iterations to decrease f_{acc}.

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Improve G for a fixed forgery detector D.

Algorithm 1 Minibatch stochastic gradient descent training of generative adversarial nets. The number of steps to apply to the discriminator, k, is a hyperparameter. We used k = 1, the least expensive option, in our experiments.

for number of training iterations do

for k steps do

- Sample minibatch of m noise samples $\{z^{(1)}, \ldots, z^{(m)}\}$ from noise prior $p_g(z)$.
- Sample minibatch of m examples $\{x^{(1)}, \dots, x^{(m)}\}$ from data generating distribution $p_{\text{data}}(x)$.
- Update the discriminator by ascending its stochastic gradient:

$$\nabla_{\theta_d} \frac{1}{m} \sum_{i=1}^m \left[\log D\left(\boldsymbol{x}^{(i)} \right) + \log \left(1 - D\left(G\left(\boldsymbol{z}^{(i)} \right) \right) \right) \right].$$

end for

- Sample minibatch of m noise samples $\{z^{(1)}, \ldots, z^{(m)}\}$ from noise prior $p_g(z)$.
- Update the generator by descending its stochastic gradient:

$$\nabla_{\theta_g} \frac{1}{m} \sum_{i=1}^{m} \log \left(1 - D\left(G\left(\boldsymbol{z}^{(i)}\right) \right) \right).$$

end for

The gradient-based updates can use any standard gradient-based learning rule. We used momentum in our experiments.

Consider the loss term for fake data:

$$f_{\text{acc}}(\theta_G, \theta_D) = \mathbb{E}_{\mathbf{x} \sim P_{\text{data}}(\mathbf{x})}[\log D_{\theta_D}(\mathbf{x})] + \mathbb{E}_{\mathbf{z} \sim P(\mathbf{z})}[\log(1 - D_{\theta_D}(G_{\theta_G}(\mathbf{z})))]$$

 What happens early during training, when G is not very good (but D typically is fairly good)?

$$\nabla_{\theta_G} \log(1 - D(G(\mathbf{z}))) = -\frac{1}{1 - D(G(\mathbf{z}))} \frac{\partial D}{\partial G} \frac{\partial G}{\partial \theta_G}$$

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$$= -\frac{1}{1} \sigma'(v) \frac{\partial v}{\partial G} \frac{\partial G}{\partial \theta_G}$$

Here we assume D uses a logistic sigmoid σ as its output layer, whose input is v.

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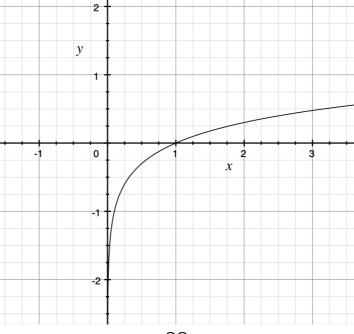
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$$= -\frac{1}{1} \sigma'(v) \frac{\partial v}{\partial G} \frac{\partial G}{\partial \theta_G}$$
$$\approx -1 \times 0 \times \frac{\partial v}{\partial G} \frac{\partial G}{\partial \theta_G}$$

 To accelerate training early on, we can instead use a different loss term for the fake data that yields the same desired behavior but trains faster.

$$f_{\text{acc}}(\theta_G, \theta_D) = \mathbb{E}_{\mathbf{x} \sim P_{\text{data}}(\mathbf{x})} [\log D_{\theta_D}(\mathbf{x})] + \mathbb{E}_{\mathbf{z} \sim P(\mathbf{z})} [-\log(D_{\theta_D}(G_{\theta_G}(\mathbf{z})))]$$

New loss term

 The reason is that the gradient of -log(v) for v≈0 is very large, whereas the gradient of log(1-v)≈1.

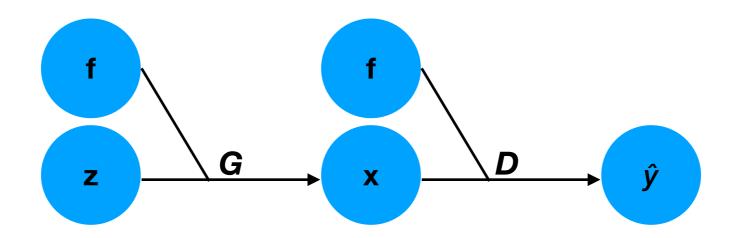


GANS

- Show Goodfellow et al. 2014 paper.
- GANs represent the state-of-the-art (as of 2021) for generating realistic data.
- GANs have also inspired many other adversarial training methods.

Conditional GANs

- One example is conditional GANs:
 - G also accepts a parameter vector f (e.g., 1-hot encoding of MNIST class) that specifies what kind of data to generate.
 - D also accepts f to help discriminate a particular kind of real from fake data.



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- GANs are renowned for being difficult to train:
 - 1. How to choose k, l? More hyperparameters to optimize.
 - 2.We will probably never reach the equilibrium where G exactly produces $P_{\text{data}}(\mathbf{x})$ and D's accuracy is 0.5.
 - What kind of "training curve" for D, G should we expect?
 - If *D* gets too good too fast, then *G* may never have a chance to improve.

- GANs are renowned for being difficult to train:
 - 3. Mode collapse G generates realistic data but only for a *subset* of the domain of $P_{\text{data}}(\mathbf{x})$, e.g.:

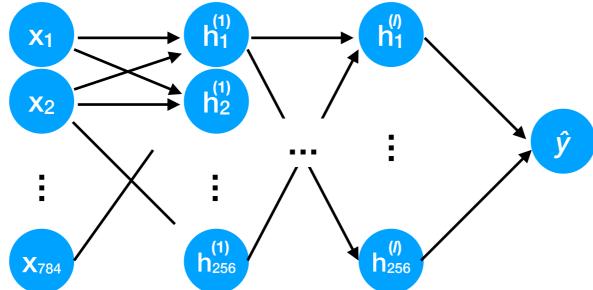


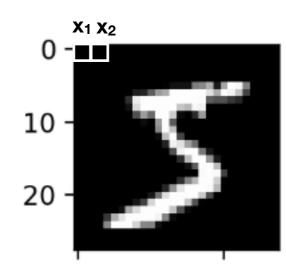
- GANs are renowned for being difficult to train:
 - 3. Neuron co-adaptation training gets stuck because multiple pathways rely on each other too much.

Consider an MNIST image near the borders: what

property do pixels x_1 , x_2 have?

•
$$x_1 = x_2 = 0$$
 ?



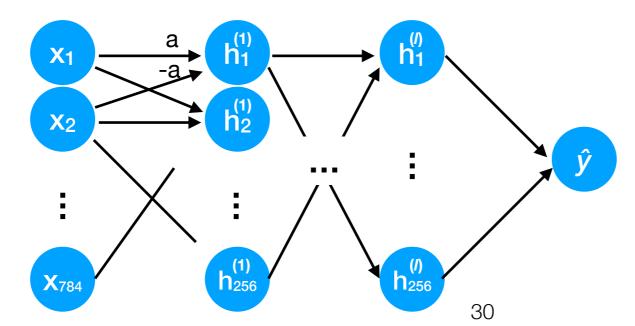


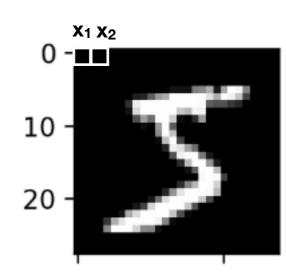
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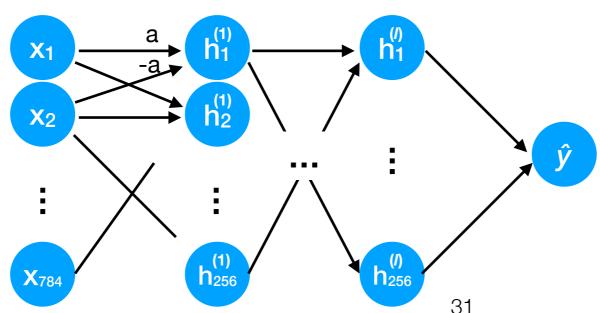




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 In the latter case, D gives feedback to G that images are "ok" as long the background noise "cancels"

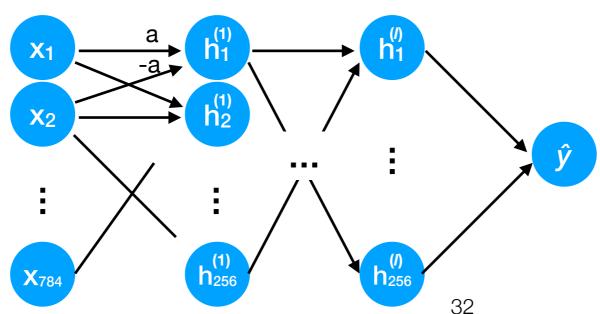
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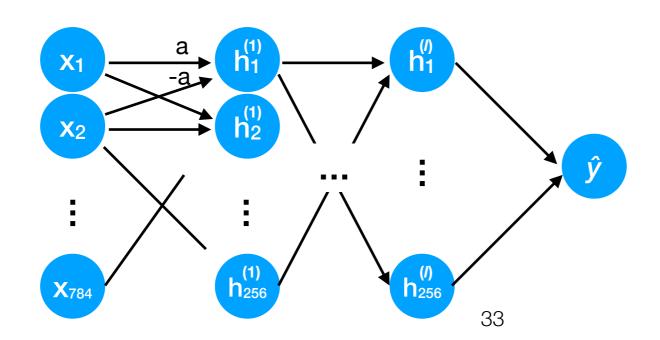
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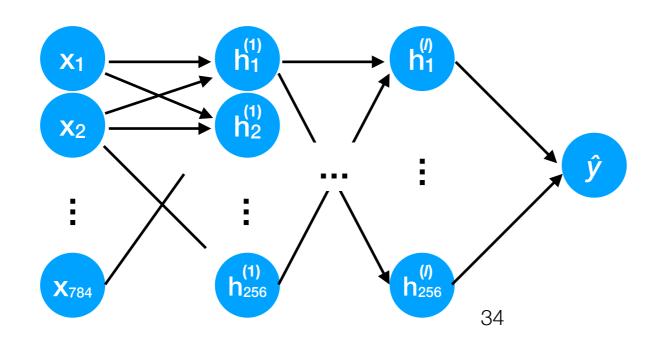
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- GANs are renowned for being difficult to train:
 - 3. Neuron co-adaptation training gets stuck because multiple pathways rely on each other too much.

• To prevent this from occurring, we can use dropout on the input layer **x**, so that each pathway must judge

independently if the image is a fake.



- One of the weaknesses of DL is that we often do not know how the models work.
- This means that models can sometimes be systematically exploited to give nonsensical outputs.
- In high-stakes scenarios such as autonomous vehicles, the results can be disastrous (e.g., misread a speed limit sign of "25km/h" for "95km/h").

- One way to systematically confuse a trained NN is to "adjust" an input \mathbf{x} so as to maximally change the output \hat{y} .
- In other words, for any loss function f (e.g., cross-entropy), compute the gradient:

$$\nabla_{\mathbf{x}} f(\mathbf{x}; \mathbf{w})$$

where w are the NN's weights.

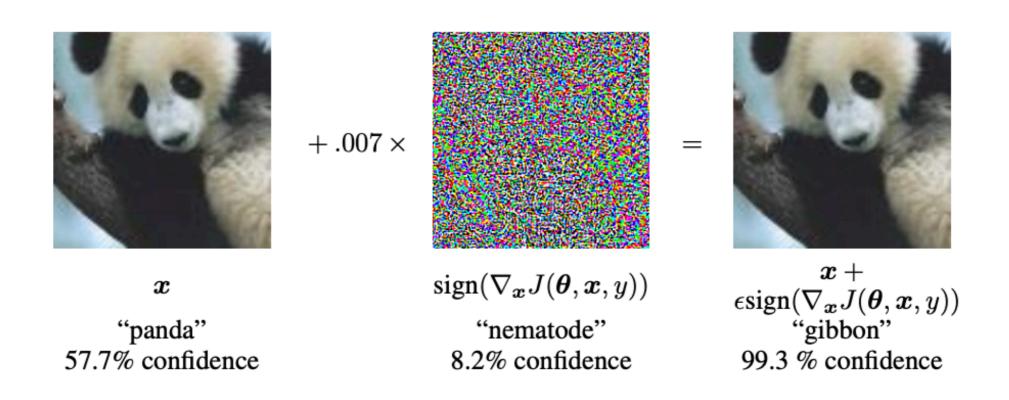
 We then construct an adversarial example (that differs by at most η from x) as:

$$\mathbf{x}' = \mathbf{x} + \eta \operatorname{sign}(\nabla_{\mathbf{x}} f(\mathbf{x}; \mathbf{w}))$$

 Note that this is different from what we usually do during training, i.e., conduct SGD on w using the gradient:

$$\nabla_{\mathbf{w}} f(\mathbf{x}; \mathbf{w})$$

 This technique is responsible for the famous panda -> gibbon mistake (Goodfellow et al. 2014):



- Since this discovery of this attack, many DL researchers have investigated how to prevent it.
- One simple technique that can help is to train explicitly on adversarial examples.
- This is easy, since we can create adversarial examples at will, and can improve robustness to outliers at test time.