Mathematics of Data: From Theory to Computation

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Lecture 11: Adversarial machine learning and generative adversarial networks

Laboratory for Information and Inference Systems (LIONS)
École Polytechnique Fédérale de Lausanne (EPFL)

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Outline

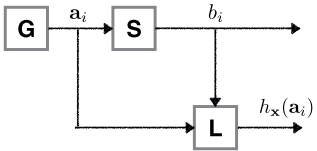
- This class
  - Adversarial Machine Learning (minmax)
    - Adversarial Training
    - Generative Adversarial Networks (GANs)
- Next class
  - Difficulty of minmax
  - Diffusion models

Adversarial machine learning

$$\min_{\mathbf{x}\in\mathcal{X}} \max_{\mathbf{y}\in\mathcal{Y}} \Phi(\mathbf{x}, \mathbf{y})$$

- A seemingly simple optimization formulation
- Critical in machine learning with many applications
  - Adversarial examples and training
  - Generative adversarial networks
  - *Robust reinforcement learning (more on this next week)
  - ...

From empirical risk minimization...



Definition (Empirical Risk Minimization (ERM))

Let $h_{\mathbf{x}} : \mathbb{R}^p \to \mathbb{R}$ be a model with parameters $\mathbf{x}$ and let $\{(\mathbf{a}_i, b_i)\}_{i=1}^n$ be samples with $b_i \in \{-1, 1\}$ and $\mathbf{a}_i \in \mathbb{R}^p$. The ERM problem reads

$$\min_{\mathbf{x}} \left\{ R_n(x) := \frac{1}{n} \sum_{i=1}^{n} L(h_{\mathbf{x}}(\mathbf{a}_i), b_i) \right\},$$

where $L(h_{\mathbf{x}}(\mathbf{a}_i), b_i)$ is the loss on the sample $(\mathbf{a}_i, b_i)$.

Some frequently used loss functions

- $L(h_{\mathbf{x}}(\mathbf{a}_i), b_i) = \log(1 + \exp(-b_i h_{\mathbf{x}}(\mathbf{a}_i)))$ *Logistic loss.*
- $L(h_{\mathbf{x}}(\mathbf{a}_i), b_i) = (b_i - h_{\mathbf{x}}(\mathbf{a}_i))^2$ *Squared error.*
- $L(h_{\mathbf{x}}(\mathbf{a}_i), b_i) = \max(0, 1 - b_i h_{\mathbf{x}}(\mathbf{a}_i))$ *Hinge loss.*

...Into adversarial examples

Definition (Adversarial examples [28])

Let $h_{\mathbf{x}^\star} : \mathbb{R}^p \to \mathbb{R}$ be a model trained through empirical risk minimization, with optimal parameters $\mathbf{x}^\star$. Let $(\mathbf{a}, b)$ be a sample with $b \in \{-1, 1\}$ and $\mathbf{a} \in \mathbb{R}^p$. An **adversarial example** is a perturbation $\boldsymbol{\delta} \in \mathbb{R}^p$ designed to lead the trained model $h_{\mathbf{x}^\star}$ to misclassify a given input $\mathbf{a}$. Given an $\epsilon > 0$, it is constructed by solving

$$\boldsymbol{\delta} \in \underset{\boldsymbol{\delta}:\|\boldsymbol{\delta}\| \leq \epsilon}{\arg\max}\, L(h_{\mathbf{x}^\star}(\mathbf{a} + \boldsymbol{\delta}), b)$$

Example norms frequently used in adversarial attacks

- The most commonly used norm is the ℓ_∞-norm [12, 21].
- The use of ℓ_1-norm leads to sparse attacks.

Figure: (Left) An ℓ_∞-attack: The alteration is hard to perceive. (Right) An ℓ_1-attack: The alteration in this case is obvious.

Adversarial examples and proximal gradient descent

- Target problem:

$$\max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|_\infty \leq \epsilon} L(h_{\mathbf{x}^\star}(\mathbf{a}+\boldsymbol{\delta}), b)$$

- We can do better than FGSM via proximal gradient methods for composite minimization:

$$\max_{\boldsymbol{\delta}\in\mathbb{R}^p} \underbrace{L(h_{\mathbf{x}^\star}(\mathbf{a}+\boldsymbol{\delta}), b)}_{f(\boldsymbol{\delta})} + \underbrace{\delta_{\mathcal{N}}(\boldsymbol{\delta})}_{g(\boldsymbol{\delta})},$$

where $\delta_{\mathcal{N}}(\boldsymbol{\delta})$ is the indicator function of the ball $\mathcal{N} := \{\boldsymbol{\delta} : \|\boldsymbol{\delta}\|_\infty \leq \epsilon\}$.

Recall: Proximal operator of indicator functions

For the indicator functions of simple sets, e.g., $g(\boldsymbol{\delta}) := \delta_{\mathcal{N}}(\boldsymbol{\delta})$, the prox-operator is the projection operator

$$\text{prox}_{\lambda g}(\boldsymbol{\delta}) := \pi_{\mathcal{N}}(\boldsymbol{\delta}),$$

where $\pi_{\mathcal{N}}(\boldsymbol{\delta})$ denotes the projection of $\boldsymbol{\delta}$ onto $\mathcal{N}$. When $\mathcal{N} = \{\boldsymbol{\delta} : \|\boldsymbol{\delta}\|_\infty \leq \lambda\}$, $\pi_{\mathcal{N}}(\boldsymbol{\delta}) = \text{clip}(\boldsymbol{\delta}, [-\lambda, \lambda])$.

Adversarial examples and proximal gradient descent (cont'd)

- Target non-convex problem:

$$\max_{\boldsymbol{\delta} \in \mathbb{R}^p} \underbrace{L(h_{\mathbf{x}^\star}(\mathbf{a}+\boldsymbol{\delta}), b)}_{f(\boldsymbol{\delta})} + \underbrace{\delta_{\mathcal{N}}(\boldsymbol{\delta})}_{g(\boldsymbol{\delta})},$$

where $\delta_{\mathcal{N}}(\boldsymbol{\delta})$ is the indicator function of the ball $\mathcal{N} := \{\mathbf{y} : \|\mathbf{y}\|_\infty \leq \epsilon\}$.

Proximal gradient ascent (PGA)

1. Choose $\boldsymbol{\delta}^0 \in \operatorname{dom} f(\boldsymbol{\delta}) + g(\boldsymbol{\delta})$ as initialization.
2. For $k = 0, 1, \cdots$, generate a sequence $\{\boldsymbol{\delta}^k\}_{k \geq 0}$ as:

$$\boldsymbol{\delta}^{k+1} := \operatorname{prox}_{\alpha_k g}\left(\boldsymbol{\delta}^k + \alpha_k \nabla f(\boldsymbol{\delta}^k)\right).$$

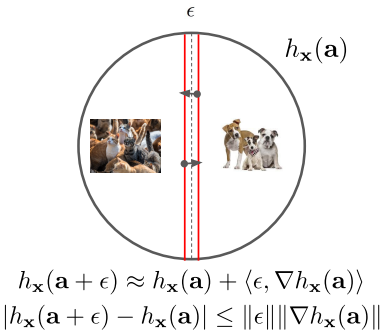
Remarks:

- PGA results in more powerful adversarial "attacks" than FGSM [16].
- The PGA is incorrectly referred to as projected gradient descent in this literature.
- Practitioners prefer to use several steps of FGSM instead of PGA [17, 18, 21]:

$$\boldsymbol{\delta}^{k+1} = \pi_{\mathcal{X}}\left(\boldsymbol{\delta}^k + \alpha_k \ \textbf{sign}\left(\nabla f(\boldsymbol{\delta}^k)\right)\right).$$

- See the appendix for a through study of the FSGM.

Challenge: Adversarial examples are inevitable



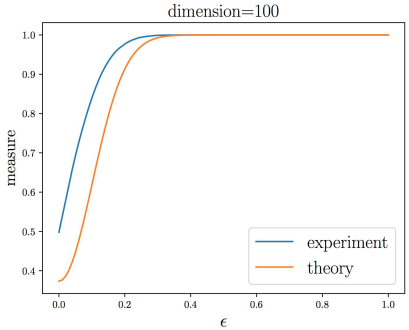
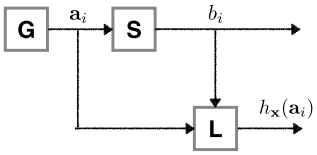


Figure: Understanding the robustness of a classifier in high-dimensional spaces. Shafahi et al. 2019.

Hardness results have never been a barrier for ML researchers



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Let $h_{\mathbf{x}} : \mathbb{R}^p \to \mathbb{R}$ be a model with parameters $\mathbf{x}$ and let $\{(\mathbf{a}_i, b_i)\}_{i=1}^n$ be samples with $b_i \in \{-1, 1\}$ and $\mathbf{a}_i \in \mathbb{R}^p$. The ERM problem reads

$$\min_{\mathbf{x}} \left\{ R_n(x) := \frac{1}{n} \sum_{i=1}^{n} L(h_{\mathbf{x}}(\mathbf{a}_i), b_i) \right\},$$

where $L(h_{\mathbf{x}}(\mathbf{a}_i), b_i)$ is the loss on the sample $(\mathbf{a}_i, b_i)$.

Objectives

- $\min_{\mathbf{x}} \left\{ \frac{1}{n} \sum_{i=1}^{n} \left[\max_{\boldsymbol{\delta}: \|\boldsymbol{\delta}\|_\infty \leq \epsilon} L\left(h_{\mathbf{x}}\left(\mathbf{a}_i + \boldsymbol{\delta}\right), b_i\right) \right] \right\}$ — Adversarial training [14].
- $\min_{\mathbf{x}} \left\{ \frac{1}{n} \sum_{i=1}^{n} \left[\max_{\boldsymbol{\delta}: \|\boldsymbol{\delta}\|_2 \leq \epsilon} L(h_{\mathbf{x}+\boldsymbol{\delta}}\left(\mathbf{a}_i\right), b_i) \right] \right\}$ — ϵ-stability training [4], Sharpness-aware minimization [10].
- $\min_{\mathbf{x}} \max_{b^c \in [C]} \frac{1}{n_c} \sum_{i=1}^{n_c} \left[\max_{\boldsymbol{\delta}: \|\boldsymbol{\delta}\| \leq \epsilon} L\left(h_{\mathbf{x}}\left(\mathbf{a}_i + \boldsymbol{\delta}\right), b^c\right) \right]$ — Class fairness [25].

Remark: ○ We focus on adversarial training during the lecture. See supplementary material for more.

Towards adversarial training

Adversarial Training [14]

Let $h_{\mathbf{x}} : \mathbb{R}^n \to \mathbb{R}$ be a model with parameters $\mathbf{x}$ and let $\{(\mathbf{a}_i, b_i)\}_{i=1}^n$, with the data $\mathbf{a}_i \in \mathbb{R}^p$ and the labels b_i. The problem of adversarial training is the following adversarial optimization problem

$$\min_{\mathbf{x}} \frac{1}{n} \sum_{i=1}^{n} \left[\max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|_\infty \leq \epsilon} L(h_{\mathbf{x}}(\mathbf{a}_i + \boldsymbol{\delta}), b_i) \right] \approx \min_{\mathbf{x}} \mathbb{E}_{(\mathbf{a},b)\sim\mathbb{P}} \left[\max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|_\infty \leq \epsilon} L(h_{\mathbf{x}}(\mathbf{a}_i + \boldsymbol{\delta}), b_i) \right].$$

Note the similarity with the template $\min_{\mathbf{x}\in\mathcal{X}} \max_{\mathbf{y}\in\mathcal{Y}} \Phi(\mathbf{x}, \mathbf{y})$.

Beyond robustness: Adversarial training for better interpretability

- Retinopathy classification problem: Given a retinal image (left), predict whether there is a disease.
- **Zeiss:** How can we interpret the prediction of a model $h_{\mathbf{x}}(\mathbf{a})$?
- **Solution:** Look at $\nabla_{\mathbf{x}} h_{\mathbf{x}}(\mathbf{a})$, called the saliency map [9]. Minimax adversarial training seems to help!

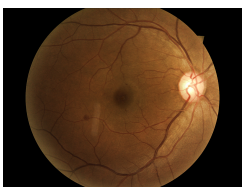
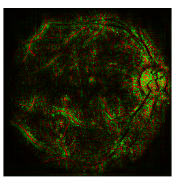
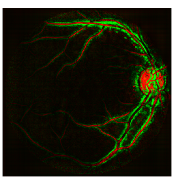

Table: **Left:** Ground truth image, **Middle:** Saliency map, **Right:** Saliency map with adversarial training.

Solving the outer problem

Adversarial Training [14]

Let $h_{\mathbf{x}} : \mathbb{R}^p \to \mathbb{R}$ be a model with parameters $\mathbf{x}$ and let $\{(\mathbf{a}_i, b_i)\}_{i=1}^n$, with $\mathbf{a}_i \in \mathbb{R}^p$ and b_i be the corresponding labels. The adversarial training optimization problem is given by

$$\min_{\mathbf{x}} \left\{ \frac{1}{n} \sum_{i=1}^{n} f_i(\mathbf{x}) := \frac{1}{n} \sum_{i=1}^{n} \underbrace{\left[\max_{\boldsymbol{\delta}: \|\boldsymbol{\delta}\|_\infty \leq \epsilon} L(h_{\mathbf{x}}(\mathbf{a}_i + \boldsymbol{\delta}), b_i) \right]}_{=: f_i(\mathbf{x})} \right\}.$$

Note that L is not continuously differentiable due to ReLU, max-pooling, etc.

Solving the outer problem

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Let $h_{\mathbf{x}} : \mathbb{R}^p \to \mathbb{R}$ be a model with parameters $\mathbf{x}$ and let $\{(\mathbf{a}_i, b_i)\}_{i=1}^n$, with $\mathbf{a}_i \in \mathbb{R}^p$ and b_i be the corresponding labels. The adversarial training optimization problem is given by

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Note that L is not continuously differentiable due to ReLU, max-pooling, etc.

Question

How can we compute the gradient

$$\nabla_{\mathbf{x}} f_i(\mathbf{x}) := \nabla_{\mathbf{x}} \left(\max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|_\infty \leq \epsilon} L(h_{\mathbf{x}}(\mathbf{a}_i + \boldsymbol{\delta}), b_i) \right)?$$

- **Challenge:** It involves differentiating with respect to a maximization.
- **A solution:** We can use Danskin's theorem under some conditions.

Danskin's Theorem (1966): How do we compute the gradient?

Theorem ([5])

Let $\mathcal{S}$ be compact set, $\Phi : \mathbb{R}^p \times \mathcal{S}$ be continuous such that $\Phi(\cdot, \mathbf{y})$ is differentiable for all $\mathbf{y} \in \mathcal{S}$, and $\nabla_{\mathbf{x}} \Phi(\mathbf{x}, \mathbf{y})$ be continuous on $\mathbb{R}^p \times \mathcal{S}$. Define

$$f(\mathbf{x}) \coloneqq \max_{\mathbf{y} \in \mathcal{S}} \Phi(\mathbf{x}, \mathbf{y}), \qquad \mathcal{S}^\star(\mathbf{x}) \coloneqq \arg\max_{\mathbf{y} \in \mathcal{S}} \Phi(\mathbf{x}, \mathbf{y}).$$

Let $\mathbf{d} \in \mathbb{R}^p$, and $\|\mathbf{d}\|_2 = 1$. The directional derivative $D_{\mathbf{d}} f(\bar{\mathbf{x}})$ of f in the direction $\mathbf{d}$ at $\bar{\mathbf{x}}$ is given by

$$D_{\mathbf{d}} f(\bar{\mathbf{x}}) = \max_{\mathbf{y} \in \mathcal{S}^\star(\bar{\mathbf{x}})} \langle \mathbf{d}, \nabla_{\mathbf{x}} \Phi(\bar{\mathbf{x}}, \mathbf{y}) \rangle.$$

An immediate consequence

If $\boldsymbol{\delta}^\star \in \arg\max_{\boldsymbol{\delta}: \|\boldsymbol{\delta}\| \leq \epsilon} L(h_{\mathbf{x}}(\mathbf{a}_i + \boldsymbol{\delta}), b_i)$ is unique, then we have

$$\nabla_{\mathbf{x}} f_i(\mathbf{x}) = \nabla_{\mathbf{x}} L(h_{\mathbf{x}}(\mathbf{a}_i + \boldsymbol{\delta}^\star), b_i).$$

A practical implementation of adversarial training: Stochastic subgradient descent

Stochastic Adversarial Training [21]

Input: learning rate α_k, iterations T, batch size K.

1. initialize neural network parameters $\mathbf{x}^0$

2. **For** $k = 0, 1, ..., T$:

i. initialize update vector $\mathbf{g}^k := 0$

ii. select a mini-batch of data $B \subset \{1, \dots, n\}$ with $|B| = K$

iii. **For** $i \in B$:

a. Find an attack $\boldsymbol{\delta}^\star$ by (approximately) solving

$$\boldsymbol{\delta}^\star \in \arg\max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|_\infty \leq \epsilon} L(h_{\mathbf{x}^k}(\mathbf{a}_i + \boldsymbol{\delta}), b_i)$$

b. Store update

$$\mathbf{g}^k := \mathbf{g}^k + \nabla_{\mathbf{x}} L(h_{\mathbf{x}^k}(\mathbf{a}_i + \boldsymbol{\delta}^\star), b_i)$$

iv. Update parameters

$$\mathbf{x}^{k+1} := \mathbf{x}^k - \frac{\alpha_k}{K}\mathbf{g}^k$$

Remarks:

- Expensive!
- Inner problem **iii.a** cannot be solved to optimality (non-convex).
- Practitioners use FGSM or PGA or PGA-ℓ_∞ to approximate the true $\boldsymbol{\delta}^\star$.
- Update in step **iii.b** is motivated by Corollary A.2 in [21]

Optimized perturbations are typically not unique!

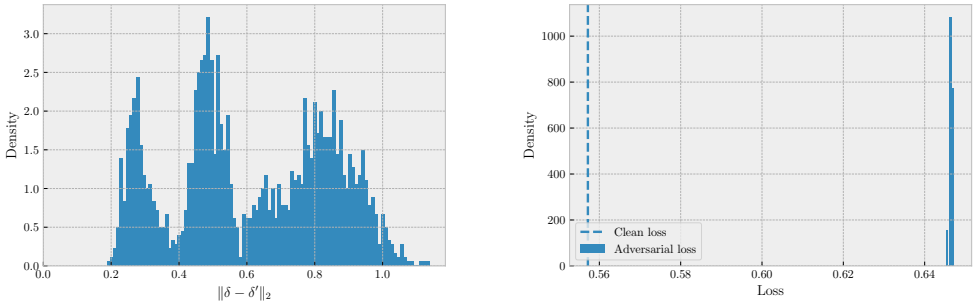

Figure: (*left*) Pairwise ℓ_2-distances between "optimized" perturbations with different initializations are bounded away from zero. (*right*) The losses of multiple perturbations on the same sample concentrate around a value much larger than the clean loss.

Theoretical foundations

	unique $\delta^\star$	non-unique $\delta^\star$
$\nabla_{\mathbf{x}}\Phi(\mathbf{x},\delta^\star)$	$\nabla_{\mathbf{x}} f(\mathbf{x})$	descent direction [21]

Published as a conference paper at ICLR 2018

Towards Deep Learning Models Resistant to Adversarial Attacks

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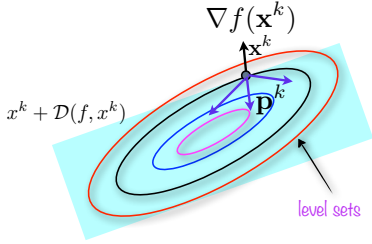

Theoretical foundations ?

	unique $\delta^\star$	non-unique $\delta^\star$
$\nabla_{\mathbf{x}}\Phi(\mathbf{x},\delta^\star)$	$\nabla_{\mathbf{x}} f(\mathbf{x})$	descent direction [21]

Published as a conference paper at ICLR 2018

TOWARDS DEEP LEARNING MODELS RESISTANT TO ADVERSARIAL ATTACKS

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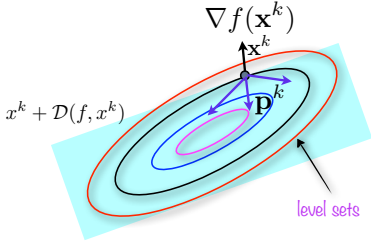
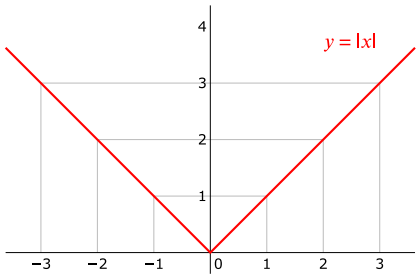


A counterexample

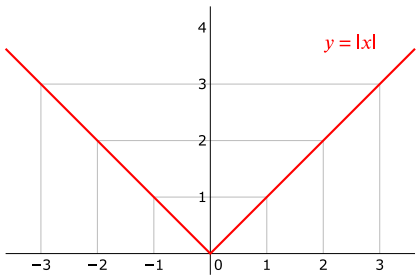
$$f(\mathbf{x}) \coloneqq \max_{\boldsymbol{\delta} \in [-1,1]} \mathbf{x}\boldsymbol{\delta} = |\mathbf{x}| \,.$$



- We have $\mathcal{S} \coloneqq [-1, 1]$ and $\Phi(\mathbf{x}, \boldsymbol{\delta}) = \mathbf{x}\boldsymbol{\delta}$.
- At $\mathbf{x} = 0$, we have $\mathcal{S}^\star(0) = [-1, 1]$.
- We can choose $\boldsymbol{\delta} = 1 \in \mathcal{S}^\star(0)$: $\Phi(\mathbf{x}, 1) = \mathbf{x}$.

A counterexample

$$f(\mathbf{x}) \coloneqq \max_{\boldsymbol{\delta} \in [-1,1]} \mathbf{x}\boldsymbol{\delta} = |\mathbf{x}| \,.$$



- We have $\mathcal{S} \coloneqq [-1, 1]$ and $\Phi(\mathbf{x}, \boldsymbol{\delta}) = \mathbf{x}\boldsymbol{\delta}$.
- At $\mathbf{x} = 0$, we have $\mathcal{S}^\star(0) = [-1, 1]$.
- We can choose $\boldsymbol{\delta} = 1 \in \mathcal{S}^\star(0)$: $\Phi(\mathbf{x}, 1) = \mathbf{x}$.
  - $-\nabla_{\mathbf{x}} \Phi(0, 1) = -1 \neq 0$.
  - Is -1 a descent direction at $\mathbf{x} = 0$?

Descent directions in the non-convex case

General Danskin's Theorem

Assume $\mathcal{Y}$ is compact and $\Phi(\mathbf{x}, \mathbf{y})$ differentiable in $\mathbf{x}$ but not necessarily convex in $\mathbf{x}$. Define $\mathcal{Y}^\star(\mathbf{x}) := \arg\max_{\mathbf{y}\in\mathcal{Y}} \Phi(\mathbf{x}, \mathbf{y})$ as the set of maximizers. Then $f(\mathbf{x}) := \max_{\mathbf{y}\in\mathcal{Y}} \Phi(\mathbf{x}, \mathbf{y})$ is *directionally differentiable* and its directional derivative is given by

$$Df(\mathbf{x}, \mathbf{d}) = \max_{\mathbf{y}^\star \in \mathcal{Y}^\star(\mathbf{x})} \langle \mathbf{d}, \nabla_{\mathbf{x}} \Phi(\mathbf{x}, \mathbf{y}^\star) \rangle \tag{1}$$

Corollary A.2 in [21] (proven wrong!)

Let $\mathbf{y}_0^\star$ be a maximizer of $\max_{\mathbf{y}\in\mathcal{Y}} \Phi(\mathbf{x}, \mathbf{y})$. Then as long as $\nabla_{\mathbf{x}} \Phi(\mathbf{x}, \mathbf{y}_0^\star)$ is non-zero, $-\nabla_{\mathbf{x}} \Phi(\mathbf{x}, \mathbf{y}_0^\star)$ is a descent direction for $f(\mathbf{x})$.

Remarks:

- The notion of directional derivative is one-sided:

$$Df(\mathbf{x}, \mathbf{d}) \coloneqq \lim_{t \to 0^+} \frac{f(\mathbf{x} + t\mathbf{d}) - f(\mathbf{x})}{t} \tag{2}$$

- Only when $\mathcal{Y}^\star(\mathbf{x}) = \{\mathbf{y}^\star\}$ is a singleton, $-\nabla_{\mathbf{x}} \Phi(\mathbf{x}, \mathbf{y}^\star)$ is *necessarily* a descent direction f.

Directional derivatives, not descent directions

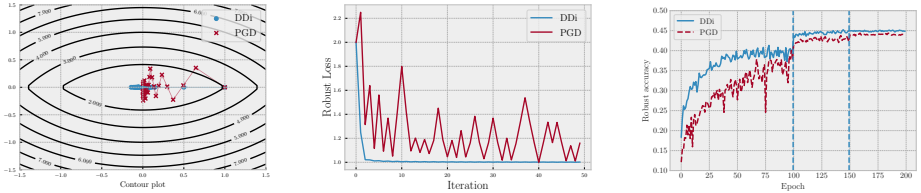


Figure: **(Left and Middle)** Synthetic adversarial training example. **(Right)** Resnet18 on CIFAR10 - Robust accuracy comparison between PGD and DDi.

Solving the inner problem does not yield a descent direction

Danskin's Theorem involves all the maximizers when computing the directional derivative along a direction $\mathbf{d}$. A single maximizer is *not* sufficient.

Remarks:

- A recent approach (DDi) computes many maximizers to find a descent direction [19].
- In practice however, the lack of descent does not seem to matter.

Danskin's theorem

Danskin's theorem (Bertsekas variant)

Let $\Phi(\mathbf{x}, \mathbf{y}) : \mathbb{R}^p \times \mathcal{Y} \to \mathbb{R}$, where $\mathcal{Y} \subset \mathbb{R}^m$ is a compact set and define $f(\mathbf{x}) := \max_{\mathbf{y} \in \mathcal{Y}} \Phi(\mathbf{x}, \mathbf{y})$. Suppose that $\Phi(\mathbf{x}, \mathbf{y})$ is convex for each $\mathbf{y}$ in the compact set $\mathcal{Y}$; the interior of the domain of f is nonempty; and $\Phi(\mathbf{x}, \mathbf{y})$ is continuous.

Define $\mathcal{Y}^\star(\mathbf{x}) := \arg\max_{\mathbf{y} \in \mathcal{Y}} \Phi(\mathbf{x}, \mathbf{y})$ as the set of maximizers and $\mathbf{y}^\star \in \mathcal{Y}^\star$ as an element of this set. We have

1. $f(\mathbf{x})$ is a convex function.
2. If $\mathcal{Y}^\star(\mathbf{x})$ is a singleton, then the function $f(\mathbf{x}) = \max_{\mathbf{y} \in \mathcal{Y}} \Phi(\mathbf{x}, \mathbf{y})$ is differentiable at $\mathbf{x}$:

$$\nabla_{\mathbf{x}} f(\mathbf{x}) = \nabla_{\mathbf{x}} \left(\max_{\mathbf{y} \in \mathcal{Y}} \phi(\mathbf{x}, \mathbf{y}) \right) = \nabla_{\mathbf{x}} \Phi(\mathbf{x}, \mathbf{y}^\star).$$

3. If $\mathcal{Y}^\star(\mathbf{x})$ contains more than one element, then the subdifferential $\partial_{\mathbf{x}} f(\mathbf{x})$ of f is given by

$$\partial_{\mathbf{x}} f(\mathbf{x}) = \operatorname{conv} \left\{ \partial_{\mathbf{x}} \Phi(\mathbf{x}, \mathbf{y}^\star) : \mathbf{y}^\star \in \mathcal{Y}^\star(\mathbf{x}) \right\}.$$

Remarks:

- The adversarial problem is not convex in $\mathbf{x}$ in general.
- (Sub)Gradients of f are calculated as $\nabla_{\mathbf{x}} f(\mathbf{x}) = \nabla_{\mathbf{x}} \Phi(\mathbf{x}, \mathbf{y}^\star)$.

Out of the frying pan into the fire

Original Formulation of Adversarial Training (I)

$$\min_{\mathbf{x}} \mathbb{E}\left[\max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|\leq\epsilon} L(h_{\mathbf{x}}(\mathbf{a}+\boldsymbol{\delta}), b)\right]$$

Original Formulation of Adversarial Training (I)

$$\min_{\mathbf{x}} \mathbb{E}\left[\max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|\leq\epsilon} L(h_{\mathbf{x}}(\mathbf{a}+\boldsymbol{\delta}), b)\right]$$

which loss L?

Original Formulation of Adversarial Training (II)

$$\min_{\mathbf{x}} \mathbb{E}\left[\max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|\leq\epsilon} L_{01}(h_{\mathbf{x}}(\mathbf{a}+\boldsymbol{\delta}), b)\right]$$

Original Formulation of Adversarial Training (II)

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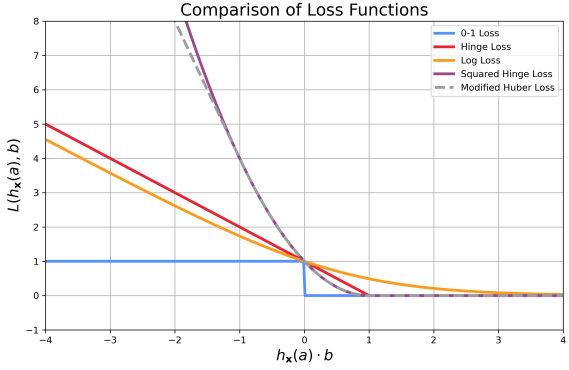
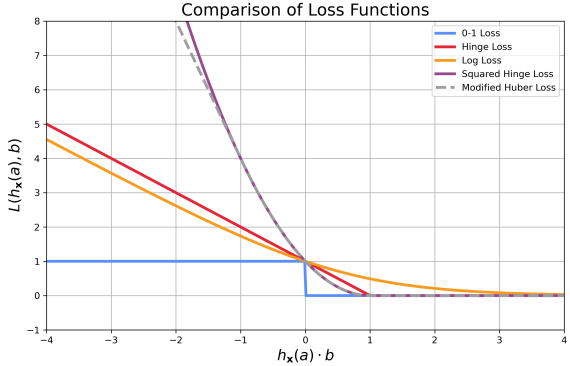
$$\min_{\mathbf{x}} \mathbb{E}\left[\max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|\leq\epsilon} L_{\mathsf{CE}}(h_{\mathbf{x}}(\mathbf{a}+\boldsymbol{\delta}), b)\right]$$

Surrogate-based optimization for Risk Minimization



Surrogate-based optimization for Risk Minimization

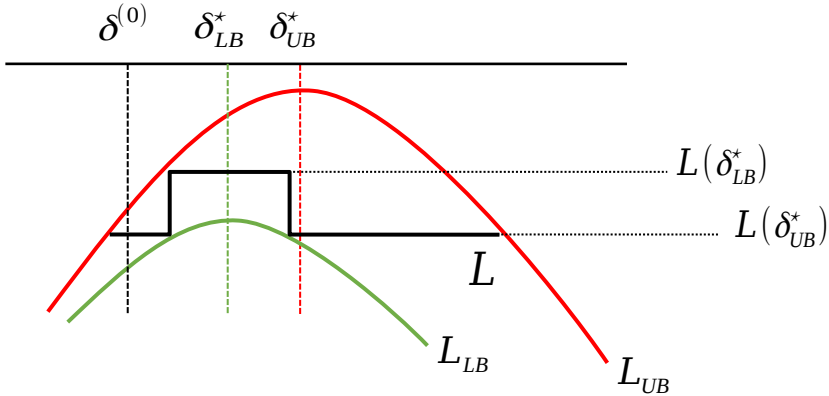


$$\mathbb{E}\left[L_{01}(h_{\mathbf{x}^\star}(\mathbf{a}+\boldsymbol{\delta}), b)\right] \leq \min_{\mathbf{x}} \mathbb{E}\left[L_{\mathsf{CE}}\left(h_{\mathbf{x}}(\mathbf{a}+\boldsymbol{\delta}), b\right)\right]$$

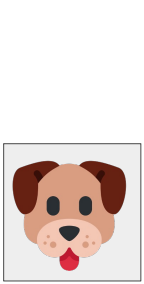
Adversary maximizes an upper bound (I)

$$L_{01}\left(h_{\mathbf{x}}(\mathbf{a}+\boldsymbol{\delta}^{\star}),b\right) \leq \max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|\leq\epsilon} L_{\mathsf{CE}}\left(h_{\mathbf{x}}(\mathbf{a}+\boldsymbol{\delta}),b\right)$$

Adversary maximizes an upper bound (II)



Why maximizing cross-entropy leads to weak adversaries



$\mathbf{a}$

$+ \quad \delta_A \quad = \quad (.49, .51, \quad 0, \quad 0)$

$L_{\mathrm{CE}}(h_{\mathbf{x}}(\mathbf{a} + \delta_A), 🐶) = 1.18$

$+ \quad \delta_B \quad = \quad (.26, .24, \quad .25, \quad .25)$

$L_{\mathrm{CE}}(h_{\mathbf{x}}(\mathbf{a} + \delta_B), 🐶) = 1.38$

Adversary's problem can be "solved" without using surrogates

Theorem (Reformulation of the Adversary's problem)

$$\boldsymbol{\delta}^\star \in \arg\max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|\leq\epsilon} \max_{j\neq b} h_{\mathbf{x}}(\mathbf{a}+\boldsymbol{\delta})_j - h_{\mathbf{x}}(\mathbf{a}+\boldsymbol{\delta})_b \Rightarrow$$

$$\boldsymbol{\delta}^\star \in \arg\max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|\leq\epsilon} L_{01}\left(h_{\mathbf{x}}(\mathbf{a}+\boldsymbol{\delta}), b\right)$$

Bilevel Optimization (BETA)

- Best targeted attack (BETA) optimization formulation [27]:

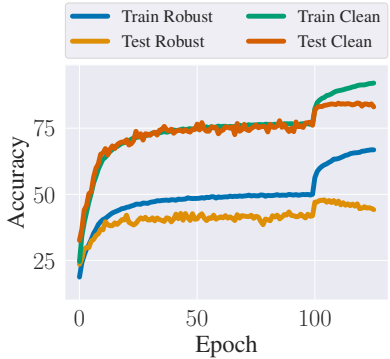
$$\min_{\mathbf{x}} \frac{1}{n} \sum_{i=1}^{n} L_{\mathsf{CE}} \left(h_{\mathbf{x}}(\mathbf{a}_i + \boldsymbol{\delta}^{\star}_{i,j^{\star}}), b_i \right)$$

$$\text{such that } \boldsymbol{\delta}^{\star}_{i,j} \in \underset{\boldsymbol{\delta}:\, \|\boldsymbol{\delta}\| \leq \epsilon}{\arg\max}\, h_{\mathbf{x}}(\mathbf{a}_i + \boldsymbol{\delta})_j - h_{\mathbf{x}}(\mathbf{a}_i + \boldsymbol{\delta})_{b_i}$$

$$j^{\star} \in \underset{j \in [K] - \{b_i\}}{\arg\max}\, h_{\mathbf{x}}(\mathbf{a}_i + \boldsymbol{\delta}_{i,j^{\star}})_j - h_{\mathbf{x}}(\mathbf{a}_i + \boldsymbol{\delta}_{i,j^{\star}})_{b_i}$$

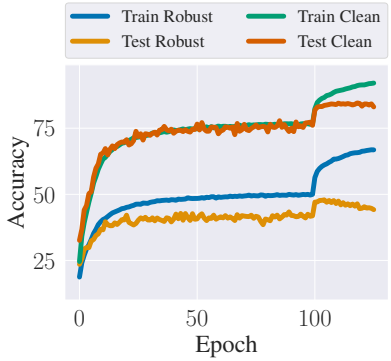
Practical Consequences of the Bilevel Formulation

Figure: Learning curves of PGD^{10}-AT (Left) and BETA^{10}-AT



Practical Consequences of the Bilevel Formulation

Figure: Learning curves of PGD^{10}-AT (Left) and BETA^{10}-AT (Right). Robust accuracy estimated with PGD^{20}



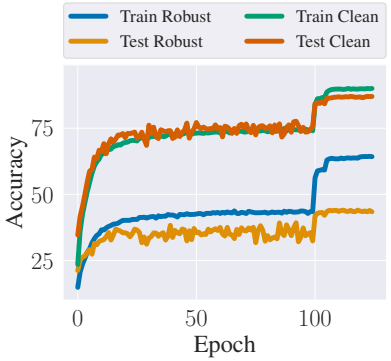


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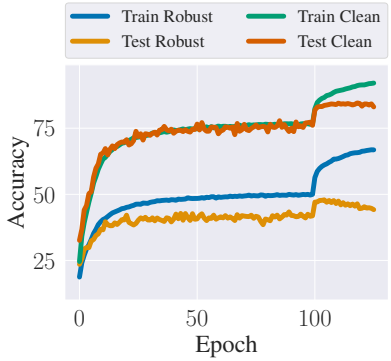
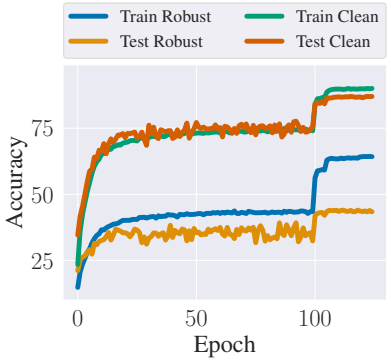
Figure: Learning curves of PGD^{10}-AT (Left) and BETA^{10}-AT (Right). Robust accuracy estimated with PGD^{20}





No Robust Overfitting occurs!

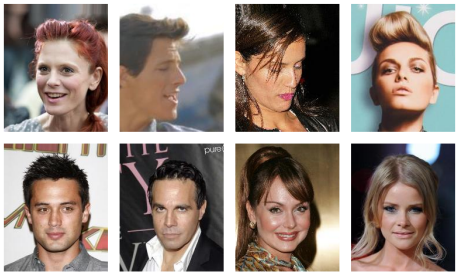
Practical Consequences of the Bilevel Formulation

Table: Adversarial performance on CIFAR-10.

Training algorithm	Test accuracy											
	Clean		FGSM		PGD^{10}		PGD^{40}		$BETA^{10}$		APGD	
	Best	Last	Best	Last	Best	Last	Best	Last	Best	Last	Best	Last
FGSM	81.96	75.43	**94.26**	**94.22**	42.64	1.49	42.66	1.62	40.30	0.04	41.56	0.00
PGD^{10}	83.71	83.21	51.98	47.39	46.74	39.90	45.91	39.45	43.64	40.21	44.36	42.62
$TRADES^{10}$	81.64	81.42	52.40	51.31	47.85	42.31	47.76	42.92	44.31	40.97	43.34	41.33
$MART^{10}$	78.80	77.20	53.84	53.73	49.08	41.12	48.41	41.55	44.81	41.22	45.00	42.90
$BETA\text{-}AT^{5}$	**87.02**	**86.67**	51.22	51.10	44.02	43.22	43.94	42.56	42.62	42.61	41.44	41.02
$BETA\text{-}AT^{10}$	85.37	85.30	51.42	51.11	45.67	45.39	45.22	45.00	44.54	44.36	44.32	44.12
$BETA\text{-}AT^{20}$	82.11	81.72	54.01	53.99	**49.96**	**48.67**	**49.20**	**48.70**	**46.91**	**45.90**	**45.27**	**45.25**

Adversarial machine learning: Introduction to Generative Adversarial Networks (GANs)

- Recall the parametric density estimation setting

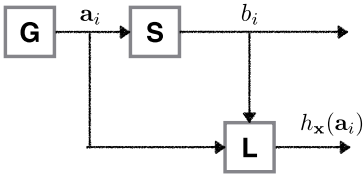


(source: http://mmlab.ie.cuhk.edu.hk/projects/CelebA.html)

$\mathbf{a}_i =$ [...images...]

$b_i =$ [...probability...]

- Goal: Games, denoising, image recovery...



- Generator $\mathbb{P}_{\mathbf{a}}$
  - Nature
- Supervisor $\mathbb{P}_{B|\mathbf{a}}$
  - Frequency data
- Learning Machine $h_{\mathbf{x}}(\mathbf{a}_i)$
  - Data scientist: Mathematics of Data

A way to model complex distributions: The push-forward measure

- Traditionally, we use analytical distributions: Restricts what we could model in real applications.
- Now, we use more expressive probability measures via *push-forward measures* with neural networks

Definition

- Let $\omega \sim \mathsf{p}_\Omega$ be a random variable.
- $h_{\mathbf{x}}(\cdot) : \mathbb{R}^p \to \mathbb{R}^m$ a function parameterized by parameters $\mathbf{x}$.

The pushforward measure of p_Ω under $h_{\mathbf{x}}$, denoted by $h_{\mathbf{x}}\#\mathsf{p}_\Omega$ is the distribution of $h_{\mathbf{x}}(\omega)$.

Example: Chi-square distribution

Let $\omega \sim \mathsf{p}_\Omega := \mathcal{N}(0,1)$ be the normal distribution. Let $h_x : \mathbb{R} \to \mathbb{R}$, $h_x(\omega) = w^x$. Let us fix $x = 2$. Then, $h_x\#\mathsf{p}_\Omega$ is the chi-square distribution with one degree of freedom.

Explanation: Change of variables.

Assume that $h : \mathbb{R}^n \to \mathbb{R}^n$ is monotonic. Given the random variable $\omega \sim \mathsf{p}_\Omega$ with probability density function $\mathsf{p}_\Omega(\omega)$, the density $\mathsf{p}_Y(\mathbf{y})$ of $\mathbf{y} = h_{\mathbf{x}}(\omega)$ reads

$$\mathsf{p}_Y(\mathbf{y}) = \mathsf{p}_\Omega(h_{\mathbf{x}}^{-1}(\mathbf{y}))\det\left(\mathbf{J}_{\mathbf{y}} h_{\mathbf{x}}^{-1}(\mathbf{y})\right)$$

where det denotes the determinant operation.

Towards an optimization problem

Problem (Ideal parametric density estimator)

Given a true distribution $\mu^\natural$, we can solve the following optimization problem,

$$\min_{\mathbf{x}} W_1(\mu^\natural, h_{\mathbf{x}} \# p_\Omega), \tag{3}$$

where the measurable function $h_{\mathbf{x}}$ is parameterized by $\mathbf{x}$ and $\omega \sim p_\Omega$ is "simple" e.g., Gaussian.

Remarks:

- See the appendix for the details of the Wasserstein distance W.
- Issues:
  - We only have access to empirical samples $\hat{\mu}_n$ of $\mu^\natural$.
  - W_1 is non-smooth, it cannot be computed exactly.

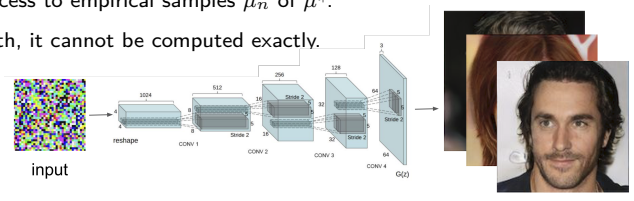
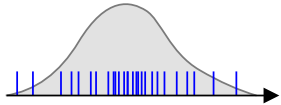


Figure: Schematic of a generative model, $h_{\mathbf{x}} \# \omega$ [11, 15].

Learning without concentration

- We can minimize $W_1\left(\hat{\mu}_n, h_{\mathbf{x}}\#\mathsf{p}_\Omega\right)$ with respect to $\mathbf{x}$.
- Figure: Empirical distribution (blue), $\hat{\mu}_n = \sum_{i=1}^n \delta_i$



A plug-in empirical estimator

Using the triangle inequality for Wasserstein distances we can upper bound in the follow way,

$$W_1(\mu^\natural, h_{\mathbf{x}}\#\mathsf{p}_\Omega) \leq W_1(\mu^\natural, \hat{\mu}_n) + W_1(\hat{\mu}_n, h_{\mathbf{x}}\#\mathsf{p}_\Omega), \tag{4}$$

where $\hat{\mu}_n$ is the empirical estimator of $\mu^\natural$ obtained from n independent samples from $\mu^\natural$.

Theorem (Slow convergence of empirical measures in 1-Wasserstein [30, 6])

Let $\mu^\natural$ be a measure defined on $\mathbb{R}^p$ and let $\hat{\mu}_n$ be its empirical measure. Then the $\hat{\mu}_n$ converges, in the worst case, at the following rate,

$$W_1(\mu^\natural, \hat{\mu}_n) \gtrsim n^{-1/p}. \tag{5}$$

Remarks:

- Using an empirical estimator in high-dimensions is terrible in the worst case.
- However, it does not directly say that $W_1\left(\mu^\natural, h_{\mathbf{x}}\#\mathsf{p}_\Omega\right)$ will be large.
- So we can still proceed and hope our parameterization interpolates harmlessly.

Duality of 1-Wasserstein

- Instead of computing W_1, we can obtain lower bounds using duality.

Theorem (Kantorovich-Rubinstein duality)

$$W_1(\mu,\nu) = \sup_{\mathtt{d}} \{\langle \mathtt{d},\mu\rangle - \langle \mathtt{d},\nu\rangle : \mathtt{d} \textit{ is 1-Lipschitz}\} \tag{6}$$

Remark:
- $\mathtt{d}$ is the "dual" variable. In the literature, it is commonly referred to as the "discriminator."

Inner product is an expectation

$$\langle \mathtt{d},\mu\rangle = \int \mathtt{d}\mathrm{d}\mu = \int \mathtt{d}(\mathbf{a})\mathrm{d}\mu(\mathbf{a}) = \boldsymbol{E}_{\mathbf{a}\sim\mu}\left[\mathtt{d}(\mathbf{a})\right]. \tag{7}$$

Kantorovich-Rubinstein duality applied to our objective

$$W_1\left(\hat{\mu}_n, h_{\mathbf{x}}\#\omega\right) = \sup\left\{\boldsymbol{E}_{\mathbf{a}\sim\hat{\mu}_n}[\mathtt{d}(\mathbf{a})] - \boldsymbol{E}_{\mathbf{a}\sim h_{\mathbf{x}}\#\omega}[\mathtt{d}(\mathbf{a})] : \mathtt{d} \text{ is 1-Lipschitz}\right\} \tag{8}$$

Integral Probability Metrics

We can define a more general class of (semi)metrics in the space of probability distributions

Definition (Integral Probability Metric)

Let $\mathcal{F}$ be a class of functions from $\mathbb{R}^p$ to $\mathbb{R}$. For two probability measures μ and ν, the IPM associated to $\mathcal{F}$ is defined as:

$$\mathcal{F}(\mu, \nu) \coloneqq \sup_{f \in \mathcal{F}} \langle f, \mu \rangle - \langle f, \nu \rangle = \sup_{f \in \mathcal{F}} \boldsymbol{E}_{\mathbf{a} \sim \mu}[f(\mathbf{a})] - \boldsymbol{E}_{\mathbf{a} \sim \nu}[f(\mathbf{a})] \tag{9}$$

Remarks:

- The 1-Wasserstein distance corresponds to $\mathcal{F} \coloneqq \{f : \mathbb{R}^p \to \mathbb{R}, f \text{ is } 1 - \text{Lipschitz}\}$
- The class cannot be described with finite parameters.

Neural network distances inspired by the 1-Wasserstein distance

- We use neural networks to parametrize a class of functions.
- Constraining the Lipschitz constant of Neural Networks is NP-Hard [29].
- We can constrain upper bounds on the Lipschitz constant [20].

Lemma

Let $h_{\mathbf{X}_1,\mathbf{X}_2}(\mathbf{a}) \coloneqq \mathbf{X}_2^T \sigma(\mathbf{X}_1\mathbf{a})$ be a one-hidden-layer neural network. Then its Lipschitz constant $L_{\mathbf{X}_1,\mathbf{X}_2}$ with respect to the ℓ_2-norm is bounded as:

$$L_{\mathbf{X}_1,\mathbf{X}_2} \leq \|\mathbf{X}_1\|_2 \|\mathbf{X}_2\|_2 \tag{10}$$

Neural Network Distance

Let

$$\mathcal{F} \coloneqq \{h_{\mathbf{X}_1,\mathbf{X}_2}(\mathbf{a}) = \mathbf{X}_2^T \sigma(\mathbf{X}_1\mathbf{a}) : \|\mathbf{X}_2\|_2 \leq 1, \|\mathbf{X}_1\|_2 \leq 1\}. \tag{11}$$

The IPM corresponding to $\mathcal{F}$ is referred to as a *Neural Network Distance*.

Remark: ○ Different network architectures/constraints lead to different Neural Network distance notions.

Wasserstein GANs formulation

- Ingredients:
  - fixed *noise* distribution p_Ω (e.g., normal)
  - target distribution $\hat{\mu}_n$ (natural images)
  - $\mathcal{X}$ parameter class inducing a class of functions (generators)
  - $\mathcal{Y}$ parameter class inducing a class of functions (dual variables)

Wasserstein GANs formulation [3]

Define a parameterized function $\mathsf{d}_{\mathbf{y}}(\mathbf{a})$, where $\mathbf{y} \in \mathcal{Y}$ such that $\mathsf{d}_{\mathbf{y}}(\mathbf{a})$ is 1-Lipschitz. In this case, the Wasserstein GAN optimization problem is given by

$$\min_{\mathbf{x}\in\mathcal{X}} \left(\max_{\mathbf{y}\in\mathcal{Y}} \boldsymbol{E}_{\mathbf{a}\sim\hat{\mu}_n} \left[\mathsf{d}_{\mathbf{y}}(\mathbf{a})\right] - \boldsymbol{E}_{\boldsymbol{\omega}\sim\mathsf{p}_\Omega} \left[\mathsf{d}_{\mathbf{y}}(h_{\mathbf{x}}(\boldsymbol{\omega}))\right] \right). \tag{12}$$

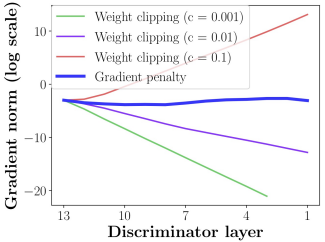
The theory-practice gap: Enforcing 1-Lipschitz of the discriminator

Weight clipping [3]

The "dual" or the "discriminator" $\mathtt{d}_{\mathbf{y}}$ weights $\mathbf{y}$ are constrained by an ℓ_∞-ball with radius $c > 0$, denoted as $\mathcal{B}$, at every iteration with

$$\pi_{\mathcal{B}}(\mathbf{y}) = \mathrm{clip}(\mathbf{y}, [-c, c]). \tag{13}$$

This trick is used to pseudo-enforce the constraint.



Remark: ○ *"Weight clipping is a clearly terrible way to enforce a Lipschitz constraint"* – original authors.

Gradient penalty [13]

Recall that 1-Lipschitz is equivalent to $\|\nabla_{\mathbf{a}} \mathtt{d}_{\mathbf{y}}(\mathbf{a})\|_* \leq 1$. This can be enforced directly through

$$\boldsymbol{E}_{\mathbf{a}\sim\hat{\mu}_n}\left[\mathtt{d}_{\mathbf{y}}(\mathbf{a})\right] - \boldsymbol{E}_{\boldsymbol{\omega}\sim\Omega}\left[\mathtt{d}_{\mathbf{y}}(h_{\mathbf{x}}(\boldsymbol{\omega}))\right] + \lambda \boldsymbol{E}_{\mathbf{a}\sim\nu}\left[\left(\|\nabla_{\mathbf{a}} \mathtt{d}_{\mathbf{y}}(\mathbf{a})\|_* - 1\right)^2\right]. \tag{14}$$

Remarks:

○ In practice the distribution ν mimicks uniform (linearly interpolated) sampling as follows: $\mathbf{a} \sim \mathrm{Uniform}(\mathbf{a}_i, h_{\mathbf{x}}(\boldsymbol{\omega}_i))$.

○ Spectral normalization: Divide each weight matrix by their spectral norm [22].

○ Learnable spline activations: both a 1-Lipschitz and more expressive architecture [24].

Practical implementation of GANs

Stochastic training of Wasserstein GANs

Input: primal and "dual" learning rates γ_t and α_m, primal iterations T, "dual" network $\mathsf{d}_{\mathbf{y}}$, generator network $h_{\mathbf{x}}$, noise distribution p_Ω, real distribution $\hat{\mu}_n$, primal and dual batch sizes B, K, "dual" iterations M.

1. initialize $\mathbf{x}^0$
2. **For** $t = 0, 1, ..., T-1$:
   - **For** $m = 0, 1, ..., M-1$:
     - initialize $\mathbf{y}^0$,
     - draw noise sample $\boldsymbol{\omega}_1, \ldots, \boldsymbol{\omega}_K \sim \mathsf{p}_\Omega$
     - draw real samples $\boldsymbol{r}_1, \ldots, \boldsymbol{r}_K \sim \hat{\mu}_n$
     - "dual" pseudo-loss $L(\mathbf{y}) := K^{-1} \sum_{i=1}^K \mathsf{d}_{\mathbf{y}}(\boldsymbol{r}_i) - \mathsf{d}_{\mathbf{y}}(h_{\mathbf{x}^t}(\boldsymbol{\omega}_i))$
     - $\sharp$update "dual" parameters $\mathbf{y}^{m+1} = \mathbf{y}^m + \gamma_m \nabla_{\mathbf{y}} L(\mathbf{y}^m)$
     - $\sharp$enforce 1-Lipschitz constraint on $\mathsf{d}_{\mathbf{y}^{m+1}}$
   - **end-For**
   - draw noise sample $\boldsymbol{\omega}_1, \ldots, \boldsymbol{\omega}_B \sim \mathsf{p}_\Omega$
   - generator pseudo-loss $L(\mathbf{x}) := -B^{-1} \sum_{i=1}^B \mathsf{d}_{\mathbf{y}^M}(h_{\mathbf{x}}(\boldsymbol{\omega}_i))$
   - update generator parameters $\mathbf{x}^{t+1} = \mathbf{x}^t - \alpha_t \nabla_{\mathbf{x}} L(\mathbf{x}^t)$
3. **end-For**

$\sharp$: Ideally, should be performed jointly.

Some historical background for a Turing award

Vanilla GAN [11]

$$\min_{\mathbf{x}\in\mathcal{X}} \max_{\mathbf{y}\in\mathcal{Y}} \boldsymbol{E}_{\mathbf{a}\sim\hat{\mu}_n} \left[\log \mathtt{d}_{\mathbf{y}}(\mathbf{a})\right] + \boldsymbol{E}_{\boldsymbol{\omega}\sim\mathsf{p}_\Omega} \left[\log\left(1 - \mathtt{d}_{\mathbf{y}}(h_{\mathbf{x}}(\boldsymbol{\omega}))\right)\right] \tag{15}$$

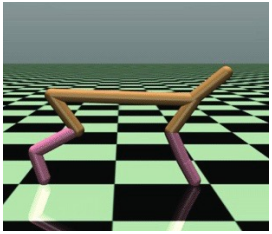
- Binary cross-entropy modeling.
- $\mathtt{d}_{\mathbf{y}}(\mathbf{a}) : \mathcal{Y} \to [0,1]$ represents the probability that $\mathbf{a}$ came from the real data distribution $\mu^\natural$.

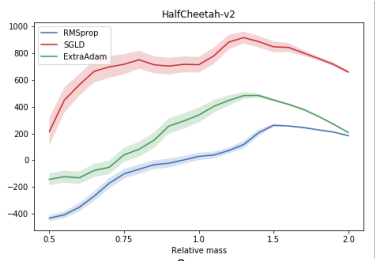
Observation: ○ Minimizes Jensen-Shannon divergence:

$$\mathrm{JSD}(\hat{\mu}_n \| h_{\mathbf{x}} \# \mathsf{p}_\Omega) = \frac{1}{2} D(\hat{\mu}_n \| h_{\mathbf{x}} \# \mathsf{p}_\Omega) + \frac{1}{2} D(h_{\mathbf{x}} \# \mathsf{p}_\Omega \| \hat{\mu}_n).$$

Take home messages

- Even the simplified view of robust & adversarial ML is challenging
- $\min$-$\max$-type has spurious attractors with no equivalent concept in $\min$-type
- Other successful attempts[1] consider "mixed Nash" concepts[2]





- Existing theory and methods for adversarial training is wrong! **... SAM too...**[3]

[1] Y-P. Hsieh, C. Liu, and V. Cevher, "Finding mixed Nash equilibria of generative adversarial networks," International Conference on Machine Learning, 2019.

[2] K. Parameswaran, Y-T. Huang, Y-P. Hsieh, P. Rolland, C. Shi, V. Cevher, "Robust Reinforcement Learning via Adversarial Training with Langevin Dynamics," NeurIPS, 2020.

[3] W. Xie, F. Latorre, K. Antonakopoulos, T. Pethick, and V. Cevher "Improving SAM requires rethinking its optimization formulation," ICLR, 2024.

Wrap up!

- Continuing on Homework 2!

A robustness example: Linear prediction

Linear model

Consider a linear model $h_{\mathbf{x}^\star}(\mathbf{a}) = \langle \mathbf{x}^\star, \mathbf{a} \rangle$ with weights $\mathbf{x}^\star \in \mathbb{R}^p$, for some input $\mathbf{a}$.

An adversarial perturbation

We aim at finding the perturbation $\boldsymbol{\delta} \in \mathbb{R}^p$ subject to $\|\boldsymbol{\delta}\|_\infty \leq \epsilon$ that produces the largest change on $h_{\mathbf{x}^\star}(\mathbf{a})$:

$$\begin{aligned}
\max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|_\infty \leq \epsilon} h_{\mathbf{x}^\star}(\mathbf{a}+\boldsymbol{\delta}) &= \max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|_\infty \leq \epsilon} \langle \mathbf{x}^\star, \mathbf{a}+\boldsymbol{\delta} \rangle \\
&= \langle \mathbf{x}^\star, \mathbf{a} \rangle + \max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|_\infty \leq \epsilon} \langle \mathbf{x}^\star, \boldsymbol{\delta} \rangle && \triangleright \text{ As } \mathbf{a} \text{ does not influence the optimization.} \\
&= \langle \mathbf{x}^\star, \mathbf{a} \rangle + \max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|_\infty \leq 1} \langle \mathbf{x}^\star, \epsilon\boldsymbol{\delta} \rangle && \triangleright \text{ By the change of variables } \boldsymbol{\delta} := \boldsymbol{\delta}/\epsilon \\
&= \langle \mathbf{x}^\star, \mathbf{a} \rangle + \epsilon\|\mathbf{x}^\star\|_1 && \triangleright \text{ Definition of the dual norm } \|\mathbf{x}\|_1 := \max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|_\infty \leq 1} \langle \mathbf{x}, \boldsymbol{\delta} \rangle
\end{aligned}$$

Taking $\boldsymbol{\delta}^\star = \operatorname{sign}(\mathbf{x}^\star)$ achieves this maximum: $\langle \mathbf{x}, \epsilon \operatorname{sign}(\mathbf{x}^\star) \rangle = \epsilon \sum_{i=1}^n \operatorname{sign}(x_i^\star) x_i^\star = \epsilon \sum_{i=1}^n |x_i^\star| = \epsilon\|\mathbf{x}^\star\|_1$.

Remarks:

- For the linear model, we have $\nabla_{\mathbf{a}} h_{\mathbf{x}^\star}(\mathbf{a}) = \mathbf{x}^\star$.
- *The gradient sign* of $h_{\mathbf{x}^\star}$ with respect to the input $\mathbf{a}$ achieves the worst perturbation.
- Sparse models are robust in linear prediction.

Adversarial examples in neural networks

- Target problem:

$$\max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|_\infty \leq \epsilon} L(h_{\mathbf{x}^\star}(\mathbf{a}+\boldsymbol{\delta}), b)$$

- Historically, researchers first tried to find approximate solutions that empirically perform well [12, 21].

Fast Gradient Sign Method (FGSM) [12]

Let $h_{\mathbf{x}^\star} : \mathbb{R}^p \to \mathbb{R}$ be a model trained through empirical risk minimization on the loss L, with optimal parameters $\mathbf{x}^\star$. Let $(\mathbf{a}, b)$ be a sample with $b \in \{-1, 1\}$ and $\mathbf{a} \in \mathbb{R}^p$. The *Fast Gradient Sign Method* computes the adversarial example

$$\boldsymbol{\delta} = \epsilon \ \text{sign}\left(\nabla_{\mathbf{a}} L(h_{\mathbf{x}^\star}(\mathbf{a}), b)\right) = \epsilon \ \text{sign}\left(\nabla_{\mathbf{a}} h_{\mathbf{x}^\star}(\mathbf{a}) \nabla_h L(h_{\mathbf{x}^\star}(\mathbf{a}), b)\right)$$

Remarks:

- The FGSM obtains adversarial examples by using *sign of the gradient of the loss*.
- Such an approach can be viewed as a linearization of the objective L around the data $\mathbf{a}$.
- For single output $h_{\mathbf{x}}(\mathbf{a})$, $\nabla_h L(h_{\mathbf{x}^\star}(\mathbf{a}), b)$ is a scalar,
  - $\text{sign}\left(\nabla_{\mathbf{a}} h_{\mathbf{x}^\star}(\mathbf{a})\right)$ pattern is important

Results of FGSM on MNIST

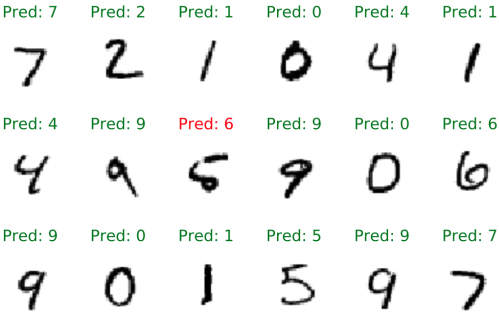


Figure: MNIST images with the predicted digit.

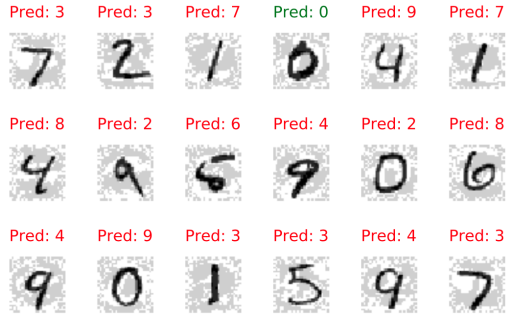


Figure: MNIST images perturbed by a FGSM attack.

Taken from `https://adversarial-ml-tutorial.org/adversarial_examples/`

A proposed link between FGSM and PGA

- Recall
  - A single step of PGA reads $\boldsymbol{\delta}_{\text{PGA}}^{k+1} := \pi_{\mathcal{N}}\left(\boldsymbol{\delta}^k + \alpha \nabla f(\boldsymbol{\delta})\right)$
  - The FGSM attack is defined as $\boldsymbol{\delta}_{\text{FGSM}} := \epsilon \ \text{sign}\left(\nabla_{\mathbf{a}} L(h_{\mathbf{x}^\star}(\mathbf{a}), b)\right)$
  - When $\mathcal{N} = \{\boldsymbol{\delta} : \|\boldsymbol{\delta}\|_\infty \leq \lambda\}$, $\pi_{\mathcal{N}}(\boldsymbol{\delta}) = \text{clip}(\boldsymbol{\delta}, [-\lambda, \lambda])$

FGSM as one step of PGA

Let $\boldsymbol{\delta}^0 = \mathbf{0}$ and $\alpha > 0$ such that $(\alpha \left|\nabla f(\mathbf{0})\right|)_i > \epsilon$ for $i = 1, \ldots, n$. Then, one step of PGA yields

$$
\begin{aligned}
\boldsymbol{\delta}_{\text{PGA}}^1 &= \pi_{\mathcal{N}}\left(\boldsymbol{\delta}^0 + \alpha \nabla_{\boldsymbol{\delta}} \nabla f(\boldsymbol{\delta}^0)\right) \\
&= \text{clip}\left(\alpha \nabla f(\mathbf{0}), [-\epsilon, \epsilon]\right) && \triangleright\ \boldsymbol{\delta}^0 = \mathbf{0} \\
&= \epsilon \ \text{sign}\left(\nabla f(\mathbf{0})\right) && \triangleright\ \text{All values are outside of the interval } [-\epsilon, \epsilon] \\
&= \epsilon \ \text{sign}\left(\nabla_{\mathbf{a}} L(h_{\mathbf{x}^\star}(\mathbf{a}), b)\right) = \boldsymbol{\delta}_{\text{FGSM}} && \triangleright\ \nabla f(\mathbf{0}) = \nabla_{\mathbf{a}} L(h_{\mathbf{x}^\star}(\mathbf{a}), b)
\end{aligned}
$$

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\end{aligned}
$$

Multiple steps of FGSM: A connection to majorization-minimization in Lecture 4

Minimization-majorization for concave functions

Let f be a concave function which is smooth in the ℓ_∞-norm with constant L_∞. Our target non-convex problem is given by

$$\max_{\boldsymbol{\delta}} f(\boldsymbol{\delta}) + \delta_{\mathcal{N}}(\boldsymbol{\delta})$$

where $\delta_{\mathcal{N}}(\boldsymbol{\delta})$ is the indicator function of the ball $\mathcal{N} := \{\boldsymbol{\delta} : \|\boldsymbol{\delta}\|_\infty \leq \epsilon\}$. Smoothness in ℓ_∞-norm implies

$$f(\boldsymbol{\delta}) + \delta_{\mathcal{N}}(\boldsymbol{\delta}) \geq \underbrace{f(\boldsymbol{\zeta}) + \langle \nabla_{\boldsymbol{\delta}} f(\boldsymbol{\zeta}), \boldsymbol{\delta} - \boldsymbol{\zeta} \rangle - \frac{L_\infty}{2} \|\boldsymbol{\delta} - \boldsymbol{\zeta}\|_\infty^2 + \delta_{\mathcal{X}}(\boldsymbol{\delta})}_{\boldsymbol{\delta}^\star \leftarrow \arg\max_{\boldsymbol{\delta}}} .$$

Maximizing the RHS with respect to $\boldsymbol{\delta}$ leads to the following (non trivial) solution [7]:

$$\boldsymbol{\delta}^\star = \text{clip}\left(\boldsymbol{\zeta} - t^\star \text{sign}(\nabla f(\boldsymbol{\zeta})), [-\epsilon, \epsilon]\right)$$

where $t^\star = \arg\max_{t:\|\boldsymbol{\delta}-\boldsymbol{\zeta}\|_\infty \leq t} \max_{\boldsymbol{\zeta}:\|\boldsymbol{\zeta}\|_\infty \leq \epsilon} \langle \nabla f(\boldsymbol{\zeta}), \boldsymbol{\delta} - \boldsymbol{\zeta} \rangle$ can be found by linear search.

Remarks:

- Setting $\boldsymbol{\zeta} = \boldsymbol{\delta}^k$ and $\boldsymbol{\delta}^\star = \boldsymbol{\delta}^{k+1}$ with a fixed step size $\alpha = t^\star$, we obtain the update in [17, 18, 21]
$$\boldsymbol{\delta}^{k+1} = \text{clip}\left(\boldsymbol{\delta}^k - t^\star \text{sign}(\nabla f(\boldsymbol{\delta}^k)), [-\epsilon, \epsilon]\right).$$
- This proof holds for **concave** and smooth functions, and need further quantification for our setting.

A notion of distance between distributions

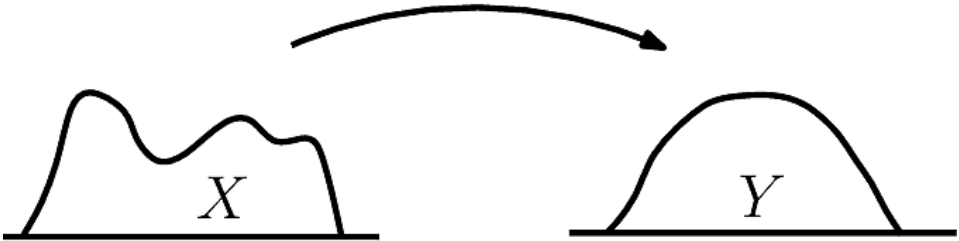

Figure: The Earth Mover's distance

Minimum cost transportation problem (Monge's problem)

Find a *transport map* $T : \mathbb{R}^d \to \mathbb{R}^d$ such that $T(X) \sim Y$, minimizing the cost

$$\text{cost}(T) := \boldsymbol{E}_X \| Y - T(X) \|. \tag{16}$$

The Wasserstein distance

Definition

Let μ and ν be two probability measures on $\mathbb{R}^d$. Their set of couplings is defined as

$$\Gamma(\mu,\nu) := \{\pi \text{ prob. measure on } \mathbb{R}^d \times \mathbb{R}^d \text{ with marginals } \mu,\nu\} \tag{17}$$

Definition (q-Wasserstein distance (Primal))

$$W_q(\mu,\nu) := \left(\inf_{\pi\in\Gamma(\mu,\nu)} \boldsymbol{E}_{(\mathbf{a},\mathbf{a}')\sim\pi} d(\mathbf{a},\mathbf{a}')^q\right)^{1/q} \tag{18}$$

where $q = 1, 2$ and d is a distance.

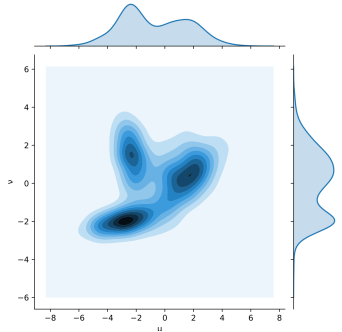

Figure: Two one-dimensional distributions plotted on the x and y axes, and one possible joint distribution that defines a transport plan between them (`https://en.wikipedia.org/wiki/Wasserstein_metric`).

Properties of the Wasserstein distance

- For any $q \geq 1$, the q-Wasserstein distance *is* a distance:
  - $W_q(\mu, \nu) = 0$ if and only if μ, ν have the same density almost everywhere (identity).
  - $W_q(\mu, \nu) = W_q(\nu, \mu)$ (symmetry).
  - $W_q(\mu, \rho) \leq W_q(\mu, \nu) + W_q(\nu, \rho)$ (triangle inequality).

Problem (Wasserstein Projection)

Given a target probability measure μ on $\mathbb{R}^d$ we are interested in solving the following optimization problem:

$$\min_{\nu \in \Delta} W_q(\mu, \nu), \tag{19}$$

where Δ is a set of probability measures on $\mathbb{R}^d$, and q is often selected as 1 or 2.

General diagram of GANs

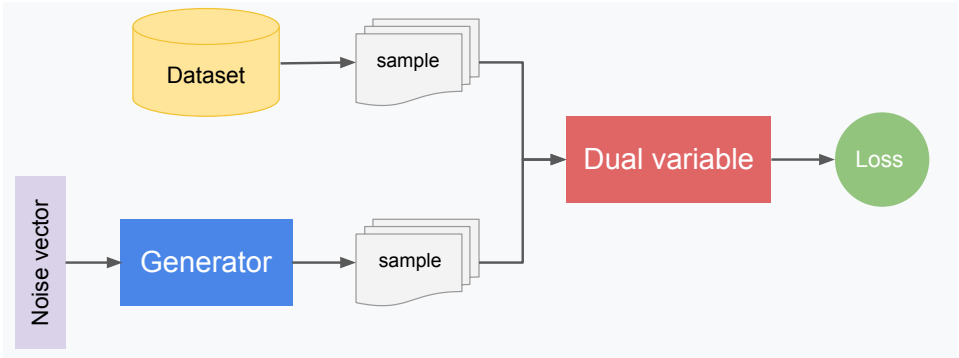

Figure: Generator/dual variable/dataset relation in GANs

*Sharpness-aware minimization (SAM) [10]

- Intuition: Flat minima usually generalizes better than sharp minima.

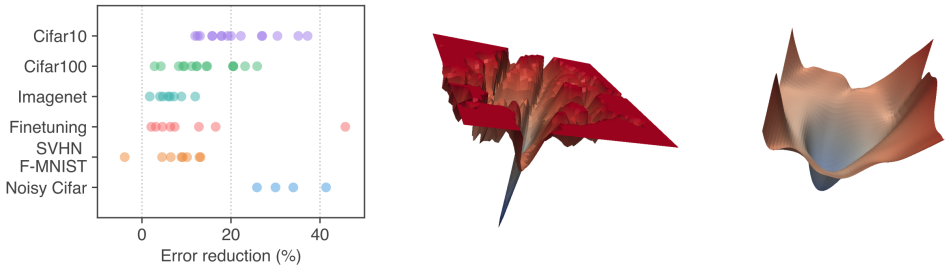

Figure: ResNet trained via SAM converges to a flatter minima **(Right)** compared with the one trained via SGD **(Middle)**, and thus leads to considerable error rate reduction **(Left)** [10].

*Sharpness-aware minimization (SAM) [10]

- Efficient approximation to the objective $\min_{\mathbf{x}} \left\{ \frac{1}{n} \sum_{i=1}^{n} \left[\max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|_2 \leq \epsilon} L(h_{\mathbf{x}+\boldsymbol{\delta}}(\mathbf{a}_i), b_i) \right] \right\}$:
  - Let's first consider the the inner maximization problem. By first-order Taylor expansion, we have:

$$\boldsymbol{\delta}^\star = \arg\max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|_2 \leq \epsilon} L\left(h_{\mathbf{x}+\boldsymbol{\delta}}(\mathbf{a}_i), b_i\right) \approx \arg\max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|_2 \leq \epsilon} \left[L\left(h_{\mathbf{x}}(\mathbf{a}_i), b_i\right) + \boldsymbol{\delta}^\top \nabla_{\mathbf{x}} L\left(h_{\mathbf{x}}(\mathbf{a}_i), b_i\right) \right]$$

$$= \arg\max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|_2 \leq \epsilon} \boldsymbol{\delta}^\top \nabla_{\mathbf{x}} L\left(h_{\mathbf{x}}(\mathbf{a}_i), b_i\right) = \epsilon \frac{\nabla_{\mathbf{x}} L\left(h_{\mathbf{x}}(\mathbf{a}_i), b_i\right)}{\|\nabla_{\mathbf{x}} L\left(h_{\mathbf{x}}(\mathbf{a}_i), b_i\right)\|_2}.$$

  - Plugging $\boldsymbol{\delta}^\star$ back the original objective and take the derivative:

$$\nabla_{\mathbf{x}} \left\{ \frac{1}{n} \sum_{i=1}^{n} \left[\max_{\boldsymbol{\delta}:\|\boldsymbol{\delta}\|_2 \leq \epsilon} L(h_{\mathbf{x}+\boldsymbol{\delta}}(\mathbf{a}_i), b_i) \right] \right\} = \frac{1}{n} \sum_{i=1}^{n} \left[\nabla_{\mathbf{x}} L(h_{\mathbf{x}+\boldsymbol{\delta}^\star}(\mathbf{a}_i), b_i) \right]$$

$$= \frac{1}{n} \sum_{i=1}^{n} \left[(1 + \frac{d\boldsymbol{\delta}^\star}{d\boldsymbol{w}}) \nabla_{\mathbf{x}} L(h_{\mathbf{x}}(\mathbf{a}_i), b_i) \mid_{\mathbf{x}+\boldsymbol{\delta}^\star} \right] \approx \frac{1}{n} \sum_{i=1}^{n} \left[\nabla_{\mathbf{x}} L(h_{\mathbf{x}}(\mathbf{a}_i), b_i) \mid_{\mathbf{x}+\boldsymbol{\delta}^\star} \right],$$

    where in the last equation the second-order term is dropped for accelerating the computation.

  - Thus, the parameters are updated by: $\mathbf{x}^{k+1} = \mathbf{x}^k - \gamma_k \frac{1}{n} \sum_{i=1}^{n} \left[\nabla_{\mathbf{x}^k} L(h_{\mathbf{x}^k}(\mathbf{a}_i), b_i) \mid_{\mathbf{x}^k + \boldsymbol{\delta}^{\star k}} \right]$, where γ_k is a step-size.

SAM's update rule

SAM

The SAM update rule is given by:

$$\tilde{\mathbf{x}}^k = \mathbf{x}^k + \epsilon \frac{\nabla L(\mathbf{x}^k)}{\|\nabla L(\mathbf{x}^k)\|} \qquad \text{[Perturb weights]}$$

$$\mathbf{x}^{k+1} = \mathbf{x}^k - \gamma_k \nabla L(\tilde{\mathbf{x}}^k) \qquad \text{[Update step]}$$

where γ_k is the step-size, and $L(\mathbf{x}) = \frac{1}{n}\sum_{i=1}^{n} L\left(h_{\mathbf{x}}\left(\mathbf{a}_i\right), b_i\right)$.

Remarks:

- SAM requires two gradient computations per update, which are sequentially dependent.
- The computational time of SAM is doubled compared with base optimizers (e.g. SGD).

Question:

- Can we run it as fast as base optimizers?

  Yes! Parallelize two gradient computations.

SAM Parallelized (SAMPa) [31]

SAMPa-λ

An auxiliary sequence $\mathbf{y}$ is introduced for parallel computation. The SAMPa update rule is given by:

$$\tilde{\mathbf{x}}^k = \mathbf{x}^k + \epsilon \frac{\nabla L(\mathbf{y}^k)}{\|\nabla L(\mathbf{y}^k)\|} \qquad \text{[Perturb weights]}$$

$$\mathbf{y}^{k+1} = \mathbf{x}^k - \gamma_k \nabla L(\mathbf{y}^k) \qquad \text{[Auxiliary sequence]}$$

$$\mathbf{x}^{k+1} = \mathbf{x}^k - (1-\lambda)\gamma_k \nabla L(\tilde{\mathbf{x}}^k) - \lambda \gamma_k \nabla L(\mathbf{y}^{k+1}) \qquad \text{[Update step]}$$

where $\lambda \in [0, 1]$. Note that $\nabla L(\tilde{\mathbf{x}}^k)$ and $\nabla L(\mathbf{y}^{k+1})$ are computed in parallel, incorporating optimistic gradient descent as follows:

$$y_{t+1} = x_t - \eta_t \nabla f(y_t), \quad x_{t+1} = x_t - \eta_t \nabla f(y_{t+1})$$

Table: **Resnet-56 with Efficient SAM variants on CIFAR-10.** The best result is in bold and the second best is underlined.

	SAM	SAMPa-0.2	LookSAM	AE-SAM	SAF	MESA	ESAM
Accuracy	94.26	**94.62**	91.42	<u>94.46</u>	93.89	94.23	94.21
Time/Epoch (s)	18.81	<u>10.94</u>	16.28	13.47	**10.09**	15.43	15.97

Fast gradient sign method (FGSM) [12]

Projected gradient descent (PGD) attack: A misnomer

Let $\boldsymbol{\eta}^{(0)} = \mathbf{0}$, the PGD update rule is given by:

$$\hat{\boldsymbol{\eta}}^{(t)} = \boldsymbol{\eta}^{(t-1)} + \alpha \cdot \operatorname{sign}\left(\nabla_{\boldsymbol{\eta}} L\left(h_{\mathbf{x}}\left(\mathbf{a} + \boldsymbol{\eta}^{(t-1)}\right), b\right)\right) \qquad \text{[Gradient step]}$$

$$\boldsymbol{\eta}^{(t)} = \max\left\{\min\left\{\hat{\boldsymbol{\eta}}^{(t)}, \epsilon\right\}, -\epsilon\right\}, \qquad \text{[Projection step]}$$

where α is the step-size and the procedure is ran for T steps. If $T = 1$ and $\alpha = \epsilon$ we recover the FGSM:

$$\boldsymbol{\eta}_{\text{FGSM}} = \epsilon \cdot \operatorname{sign}\left(\nabla_{\boldsymbol{\eta}} L\left(h_{\mathbf{x}}\left(\mathbf{a}\right), b\right)\right)$$

Problems:

- In Adversarial Training: $\times T$ overhead in training time.
- If $T = 1$, we can observe **Catastrophic Overfitting (CO)**

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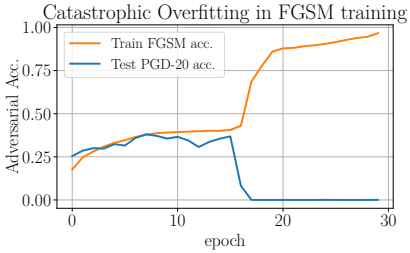
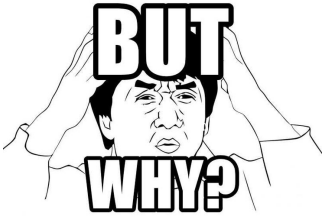
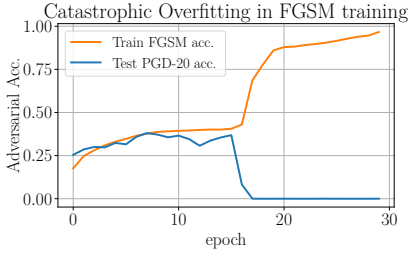
Example:

- PreActResNet18 on CIFAR10 at $\epsilon = 8/255$.
- Outcome: 100% robust to FGSM attacks and 0% robust to PGD-20 attacks.



More on CO





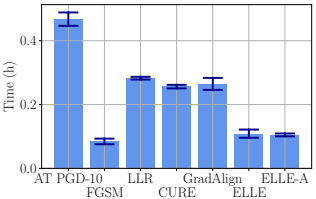
Why?

The single step solution $\boldsymbol{\eta}_{\text{FGSM}}$ only makes sense if our loss is locally linear, i.e.:

$$L\left(h_{\mathbf{x}^k}(\mathbf{a}+\boldsymbol{\eta}),b\right) \approx L\left(h_{\mathbf{x}^k}(\mathbf{a}),b\right) + \boldsymbol{\eta}^\top \nabla_{\mathbf{a}} L\left(h_{\mathbf{x}^k}(\mathbf{a}),b\right)\ , \quad \forall \boldsymbol{\eta} : ||\boldsymbol{\eta}||_\infty \leq \epsilon\,. \quad \text{[1st order Taylor expansion]}$$

Observation: This property is lost during AT with FGSM and CO appears [2].

The ELLE way [Abad Rocamora, Liu, Chrysos, Olmos and Cevher, ICLR 2024]



(a) Training time comparison

ϵ	8		16	
Method	AutoAttack	Clean	AutoAttack	Clean
LLR	$42.18 \pm (0.20)$	$75.02 \pm (0.09)$	$16.92 \pm (0.20)$	$42.81 \pm (9.62)$
CURE	$43.60 \pm (0.17)$	$77.74 \pm (0.11)$	$\underline{18.25} \pm (0.45)$	$52.49 \pm (0.04)$
GradAlign	$\mathbf{44.66} \pm (0.21)$	$\mathbf{80.50} \pm (0.07)$	$17.46 \pm (1.71)$	$44.35 \pm (15.32)$
ELLE	$42.78 \pm (0.95)$	$\underline{80.13} \pm (0.32)$	$\mathbf{18.28} \pm (0.17)$	$\mathbf{59.73} \pm (0.16)$
ELLE-A	$\underline{44.32} \pm (0.04)$	$79.81 \pm (0.10)$	$18.03 \pm (0.15)$	$\underline{59.21} \pm (1.23)$
AT PGD-10	$46.95 \pm (0.11)$	$79.11 \pm (0.08)$	$24.77 \pm (0.26)$	$59.64 \pm (0.46)$

(b) PreActResNet18 in CIFAR10

Algorithmic approaches:

- Local linearization (LLR) [26]
- Curvature regularization (CURE) [23]
- Gradient alignment (GradAlign) [2]
- Efficient local linearity regularization (ELLE) [1]

Overcoming CO with local linearity regularization [2, 1]

Why?

The single step solution $\boldsymbol{\eta}_{\text{FGSM}}$ only makes sense if our loss is locally linear, i.e.:

$$L\left(h_{\mathbf{x}^k}(\mathbf{a}+\boldsymbol{\eta}), b\right) \approx L\left(h_{\mathbf{x}^k}(\mathbf{a}), b\right) + \boldsymbol{\eta}^\top \nabla_{\mathbf{a}} L\left(h_{\mathbf{x}^k}(\mathbf{a}), b\right) \ , \quad \forall \boldsymbol{\eta} : ||\boldsymbol{\eta}||_\infty \leq \epsilon \, . \quad \text{[1st order Taylor expansion]}$$

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Observation: This property is lost during AT with FGSM and CO appears [2].

- We can measure the how locally linear a model is with the gradient missalignment.

Gradient Missalignment [2]

Let the point $\tilde{\mathbf{a}}$ be sampled uniformly such that $||\mathbf{a} - \tilde{\mathbf{a}}||_\infty \leq \epsilon$ and the gradients $\mathbf{g} = \nabla_{\mathbf{a}} L\left(h_{\mathbf{x}^k}(\mathbf{a}), b\right)$ and $\tilde{\mathbf{g}} = \nabla_{\tilde{\mathbf{a}}} L\left(h_{\mathbf{x}^k}(\tilde{\mathbf{a}}), b\right)$. The gradient missalignment is defined as:

$$\text{Grad.Miss.}(\mathbf{x}^k, \mathbf{a}) = 1 - \frac{\mathbf{g}^\top \tilde{\mathbf{g}}}{||\mathbf{g}||_2 ||\tilde{\mathbf{g}}||_2} , \tag{20}$$

with a locally linear model at $\mathbf{a}$ having $\text{Grad.Miss.}(\mathbf{x}^k, \mathbf{a}) = 0$.

Overcoming CO with local linearity regularization [2, 1]

- **Observation:** We can regularize the Gradient Missalignment during training to avoid CO, i.e., GradAlign [2]:

$$\min_{\mathbf{x}} \frac{1}{n} \sum_{i=1}^{n} L\left(h_{\mathbf{x}}\left(\mathbf{a}_i + \boldsymbol{\eta}_{\mathsf{FGSM}}\right), b_i\right) + \lambda \cdot \mathsf{Grad.Miss.}(\mathbf{x}, \mathbf{a}_i)\,.$$

- **Remark:** differentiating $\nabla_{\mathbf{x}}\mathsf{Grad.Miss.}(\mathbf{x}, \mathbf{a}_i)$ is an expensive operation due to *Double Backpropagation* [8].
- **Question:** Can we do better?.

Overcoming CO with local linearity regularization [2, 1]

- **Observation:** We can regularize the Gradient Missalignment during training to avoid CO, i.e., GradAlign [2]:

$$\min_{\mathbf{x}} \frac{1}{n} \sum_{i=1}^{n} L\left(h_{\mathbf{x}}\left(\mathbf{a}_i + \boldsymbol{\eta}_{\mathsf{FGSM}}\right), b_i\right) + \lambda \cdot \mathsf{Grad.Miss.}(\mathbf{x}, \mathbf{a}_i)\,.$$

- **Remark:** differentiating $\nabla_{\mathbf{x}}\mathsf{Grad.Miss.}(\mathbf{x}, \mathbf{a}_i)$ is an expensive operation due to *Double Backpropagation* [8].
- **Question:** Can we do better?. **Yes!**

ELLE [1]

Let the point $\tilde{\mathbf{a}}$ be sampled uniformly such that $||\mathbf{a} - \tilde{\mathbf{a}}||_\infty \leq \epsilon$ and $\hat{\mathbf{a}} = \alpha \cdot \mathbf{a} + (1-\alpha) \cdot \tilde{\mathbf{a}}$ with α sampled uniformly from $[0, 1]$. The ELLE regularization term is defined as:

$$\mathtt{ELLE}(\mathbf{x}^k, \mathbf{a}) = \left(L\left(h_{\mathbf{x}^k}(\hat{\mathbf{a}}), b\right) - \alpha \cdot L\left(h_{\mathbf{x}^k}(\mathbf{a}), b\right) - (1-\alpha) \cdot L\left(h_{\mathbf{x}^k}(\tilde{\mathbf{a}}), b\right)\right)^2\,, \tag{21}$$

with a locally linear model at $\mathbf{a}$ having $\mathtt{ELLE}(\mathbf{x}^k, \mathbf{a}) = 0$.

- **Advantage:** Regularizing ELLE does not involve *Double Backpropagation* and can as well overcome CO [1].

Is the training “fair”?

- Another grand challenge in ML: Fairness & bias
- A concrete example: Adversarial training may sacrifice subset of classes in favor of consensus
  - CIFAR10: 51% average robust accuracy while the worst class is 23.5%
  - CIFAR100: the worst class has *zero* accuracy while the best has 76%

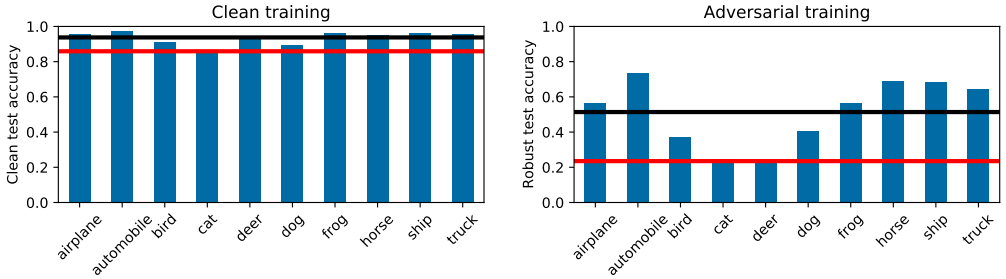


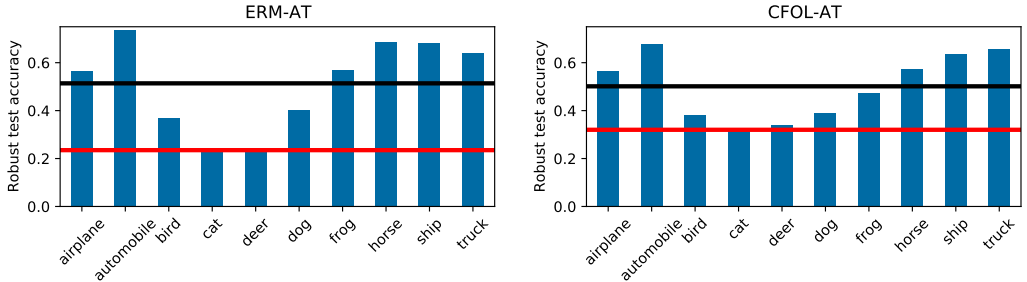
Figure: Clean accuracy and robust accuracy on CIFAR10 after clean training and adversarial training respectively.

Key challenges in ML demand much more than ERM

- Protect the weak: Class-focused online learning for adversarial training [25]

$$\min_{\mathbf{x}} \max_{b^c \in [C]} \frac{1}{n_c} \sum_{i=1}^{n_c} \left[\max_{\boldsymbol{\delta}: \|\boldsymbol{\delta}\| \leq \epsilon} L(h_{\mathbf{x}}(\mathbf{a}_i + \boldsymbol{\delta}), b^c) \right]$$

- Great potential via the minimax formulation: the average does not suffer much or can even improve!



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