# Presentation for *Probably Approximately Metric-Fair Learning*, Extension and Beyond

Probably Approximately Metric-Fair Learning is a paper by Rothblum, G. N., & Yona, G. (2018)

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#### Overview: How the Story Begins









#### **Problem**

Strict Individual Fairness does not Generalize

#### Solution

Approximate Individual Fairness Using Relaxed Metric

#### Generalization

- Bound the Loss Using Rademacher Complexity

Define
Relaxed
PACF
Learning.
Apply it to
Linear
Regression
and Logistic
Regression.

#### Thinking

This paper is theoretical work. Are there any practical ways to individual fairness and together in machine

#### Extension

Metric

Learning
Combined
with Revised
Siamese
Network for
Optimizing
both Utility
and
Individual
Fairness

### 1.Individual Fairness

#### Definition

Suppose we are given:

- 1. An input space  $\mathcal{X} = \{x_1, x_2, ..., x_n\}$ ;
- 2. A metric  $\mathcal{D}$  that defines the similarity between pair of input instances (i.e.  $\mathcal{D}(x_1, x_2)$ );
- 3. A classifier h for prediction tasks;
- 4. A distance metric d that measures the distance between output pairs (e.g.  $d(h(x_1), h(x_2))$ ).

We say the classifier h is individually fair if and only if  $\forall x_i, x_j \in \mathcal{X}, d(h(x_i), h(x_j)) \leq \mathcal{D}(x_i, x_j)$ 

Basically, the individual fairness requires the output pairs have distances strictly less than or equal to those of the according inputs.

Dwork, C., Hardt, M., Pitassi, T., Reingold, O., & Zemel, R. (2012, January)



# 2.Difficulty (Motivation for Approximation): Strict Individual Fairness is Hard to Generalize

- Trivial classifiers, i.e. generates the same outputs for all the input instances, always satisfies individual fairness both on the sample but under the entire distribution, but no utility in prediction.
- Suppose the sample is not representative of the underlying data distribution (likely to happen) or some certain representation not included in the sample, the learned classifier won't generalize to the whole distribution in terms of fairness.

# 3. Approximate Individual Fairness Metric

# Definition (loss)

The relaxed fairness loss given a similarity metric  $\Delta$  and relaxation  $\gamma$  is defined to be:

$$\ell_{\gamma,\Delta}(h,(x_1,x_2)) = \begin{cases} 1, & \text{for } d(h(x_1),h(x_2)) > \Delta(x_1,x_2) + \gamma \\ 0, & \text{for } d(h(x_1),h(x_2)) \leq \Delta(x_1,x_2) + \gamma \end{cases}$$

in which way, the individual fairness loss for the entire distribution is:

$$\mathcal{L}_{\mathcal{D},\gamma}^{F}(h) = \mathbb{E}_{x_1,x_2 \sim \mathcal{D}}[\ell_{\gamma,\Delta}(h,(x_1,x_2))]$$

the empirical loss under sample S is:

$$\hat{\mathcal{L}}_{\mathcal{D},\gamma}^{F}(h) = \frac{2}{m-1} \sum_{x_1 \in S} \sum_{x_2 \in S} [\ell_{\gamma,\Delta}(h,(x_1,x_2))]$$

# 4.PAC Learning

- Generalization error:  $\mathcal{L}(h) = \mathbb{E}_{x \sim \mathbb{D}}[\mathbb{1}(h(x) = c(x))]$  (c is the target concept, h belongs to hypothesis space  $\mathcal{H}$ )
- Empirical error:  $\hat{\mathcal{L}}_S(h) = \frac{1}{m} \sum_{x \in S} \mathbb{1}(h(x) = c(x))$

# Definition (PAC Learning)

A concept class  $\mathcal C$  is said to be PAC-learnable if  $\exists$  algorithm  $\mathcal A$  and a polynomial function  $\operatorname{poly}(,\,,\,)$  s.t. for any  $\epsilon>0$  and  $\delta>0$ , for all distributions  $\mathcal D$  on input space  $\mathcal X$  and for any target concept  $c\in\mathcal C$ , if  $m\geq \operatorname{poly}(\frac{1}{\epsilon},\frac{1}{\delta},n,\operatorname{size}(c))$ , there is:

$$\mathbb{P}_{S \sim \mathcal{D}^m}[\hat{\mathcal{L}}_S(h) \leq \epsilon] \geq 1 - \delta$$

 ${\cal C}$  is said to be efficiently PAC-learnable if  ${\cal A}$  runs in poly  $(\frac{1}{\epsilon},\frac{1}{\delta},n,{\rm size}(c))$ 

Mohri, M., Rostamizadeh, A., & Talwalkar, A. (2018)



# 5. Rademacher Complexity

# Definition (Empirical Rademacher Complexity)

For a sample  $S = \{z_1, z_2, ..., z_m\}$ , the empirical Rademacher Complexity on the hypothesis space  $\mathcal{H}$  is defined to be:

$$\hat{\mathcal{R}}_{\mathcal{S}}(\mathcal{H}) = \mathbb{E}_{\boldsymbol{\sigma}} \left[ \sup_{h \in \mathcal{H}} \frac{1}{m} \sum_{i=1}^{m} \sigma_{i} h(z_{i}) \right] = \mathbb{E}_{\boldsymbol{\sigma}} \left[ \sup_{h \in \mathcal{H}} \frac{\boldsymbol{\sigma} \cdot h(\mathbf{z})}{m} \right]$$

where the general Rademacher Complexity on the distribution  $\mathcal{D}$  is the expectation of the above one over the entire distribution:

 $\mathcal{R}_m(\mathcal{H}) = \mathbb{E}_{S \sim \mathcal{D}^m} \left[ \hat{\mathcal{R}}_S(\mathcal{H}) \right]$ . The Rademacher complexity measures to what extent the classifiers in the hypothesis space correlate with the random noise  $\sigma$ . Since it takes the supremum over the space, the more complex the space is, the more likely that it contains classifiers that correlate well with the noise.

### 6. Generalization Bounds

In order to show that this approximation generalizes well to the underlying distribution, Rothblum, G. N., & Yona, G. (2018) uses Rademacher complexity to prove a convergence bound for the loss:

#### **Theorem**

Let  $\mathcal{H}$  be a hypothesis space with Rademacher complexity  $R_m(\mathcal{H})$ .  $\forall \delta, \gamma \in (0,1), \forall G \geq 1, \forall m \geq 0 \ (w.o.l.g)$  assume m is odd, with probability at least  $1-\delta$  over an i.i.d sample  $S \sim \mathcal{D}^m$ ,  $\forall h \in \mathcal{H}$ :

$$\mathcal{M} \leq \hat{\mathcal{L}}_{\gamma}^{F}(h) \leq \mathcal{N}$$

where 
$$\mathcal{M} = \mathcal{L}_{\gamma + \frac{1}{G}}^{F}(h) - 2G(4\hat{R}_{\frac{m-1}{2}}(\mathcal{H}) + \frac{4+17\sqrt{\ln(4/\delta)}}{\sqrt{m-1}})$$
 and  $\mathcal{N} = \mathcal{L}_{\gamma + \frac{1}{G}}^{F}(h) + 2G(4\hat{R}_{\frac{m-1}{2}}(\mathcal{H}) + \frac{4+17\sqrt{\ln(4/\delta)}}{\sqrt{m-1}})$ 

# 7. How to Understand the Generalization Bounds

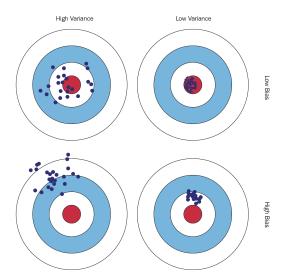
Based on the above bound, there are two essential conditions for a good generalizability:

- ullet The hypothesis space  ${\cal H}$  has a small Rademacher Complexity on the sample;
- The training sample size m is large enough.

Why should m be large? My understanding:

- 1. When there is little data in the sample, the learner is likely to choose a classifier that "looks good" based on the performance on little data. But it potentially overfits (low bias and high variance). Intuitively speaking, the learner does not have enough evidence to verify its choice.
- 2. When there is enough data in the sample (large m), the learner has enough evidence to verify its choice.

#### 8. Visualization



https://www.oreilly.com/library/view/hands-on-transfer-learning/9781788831307/1b3d8196-0617-4e1d-aa79-b2b5ef6f4a8b.xhtml

# 9.PACF Learnability

#### **Definition**

A hypothesis space  $\mathcal{H}$  is PACF learnable w.r.t a univariate accuracy loss function  $\ell^U: \mathcal{H} \times Z \to \mathbb{R}_+$  and a bivariate fairness loss function  $\ell^F_d: \mathcal{H} \times Z \times Z \to \mathbb{R}_+$  if there exists:

- a function  $m_{\mathcal{H}}(\alpha, \gamma, \delta, \epsilon, \epsilon_{\alpha}, \epsilon_{\gamma}) : (0, 1)^{6} \to \mathbb{N}$  (polynomial in  $\alpha, \gamma, \log \frac{1}{\delta}, \epsilon, \epsilon_{\alpha}, \epsilon_{\gamma}$ )
- a learning algorithm with the following property: For every required fairness parameters  $\alpha, \gamma \in (0,1)$ , failure probability  $\delta \in (0,1)$  and error parameters  $\epsilon, \epsilon_{\alpha}, \epsilon_{\gamma} \in (0,1)$  and for every distribution  $\mathcal{D}$  over  $Z = \mathcal{X} \times \mathcal{Y}$ , when running the learning algorithm on  $m \geq m_{\mathcal{H}}(\alpha, \gamma, \delta, \epsilon, \epsilon_{\alpha}, \epsilon_{\gamma})$  i.i.d examples generated by  $\mathcal{D}$ , the algorithm returns a hypothesis h such that with probability at least  $1 \delta$ :
  - $\mathcal{L}_{\mathcal{D},\gamma,d}^{Fairness}(h) \leq \alpha$ , where  $\mathcal{L}_{\mathcal{D}}^{Fairness}(h) = \mathbb{E}_{z,z' \sim \mathcal{D}}[\ell_{\gamma}^F(h,z,z')]$
  - $\mathcal{L}_{\mathcal{D}}^{Accuracy}(h) \leq \mathcal{L}_{\mathcal{D}}^{Accuracy}(\mathcal{H}^{\alpha-\epsilon_{\alpha},\gamma-\epsilon_{\gamma}}) + \epsilon$ , where  $\mathcal{L}_{\mathcal{D}}^{Accuracy}(h) = \mathbb{E}_{z \sim \mathcal{D}}[\ell_{\gamma}^{A}(h,z)]$



# 10. Relaxed PACF Learnability

# Definition $(g(\cdot)$ -relaxed PACF learnability)

A hypothesis space  $\mathcal{H}$  is  $g(\cdot)$ -relaxed PACF learnable w.r.t a function  $g:[0,1]^2 \to [0,1]^2$ , a univariate accuracy loss function  $\ell^U:\mathcal{H}\times Z\to \mathbb{R}_+$  and a bivariate fairness loss function  $\ell^F_d:\mathcal{H}\times Z\times Z\to \mathbb{R}_+$  if there exists:

- a function  $m_{\mathcal{H}}(\alpha, \gamma, \delta, \epsilon, \epsilon_{\alpha}, \epsilon_{\gamma}) : (0, 1)^{6} \to \mathbb{N}$  (polynomial in  $\alpha, \gamma, \log \frac{1}{\delta}, \epsilon, \epsilon_{\alpha}, \epsilon_{\gamma}$ )
- a learning algorithm with the following property: For every required fairness parameters  $\alpha, \gamma \in (0,1)$ , failure probability  $\delta \in (0,1)$  and error parameters  $\epsilon, \epsilon_{\alpha}, \epsilon_{\gamma} \in (0,1)$  and for every distribution  $\mathcal{D}$  over  $Z = \mathcal{X} \times \mathcal{Y}$ , when running the learning algorithm on  $m \geq m_{\mathcal{H}}(\alpha, \gamma, \delta, \epsilon, \epsilon_{\alpha}, \epsilon_{\gamma})$  i.i.d examples generated by  $\mathcal{D}$ , the algorithm returns a hypothesis h such that with probability at least  $1 \delta$ :
  - $\bullet \ \ \mathcal{L}^{\textit{Fairness}}_{\mathcal{D},\gamma,d}(\textit{h}) \leq \alpha \textit{, where } \mathcal{L}^{\textit{Fairness}}_{\mathcal{D}}(\textit{h}) = \mathbb{E}_{\textit{z},\textit{z}' \sim \mathcal{D}}[\ell^{\textit{F}}_{\gamma}(\textit{h},\textit{z},\textit{z}')]$
  - $\mathcal{L}_{\mathcal{D}}^{Accuracy}(h) \leq \mathcal{L}_{\mathcal{D}}^{Accuracy}(\mathcal{H}^{g(\alpha,\gamma)-(\epsilon_{\alpha},\epsilon_{\gamma})}) + \epsilon$ , where  $\mathcal{L}_{\mathcal{D}}^{Accuracy}(h) = \mathbb{E}_{z \sim \mathcal{D}}[\ell_{\gamma}^{A}(h,z)]$

# 11.PACF Learnability for Linear/Logistic Regression

Now it is time to apply relaxed PACF learnability into linear and logistic classifiers.

1. In linear settings,  $H_{lin} \stackrel{def}{=} \left\{ \mathbf{x} 
ightarrow rac{1 + \left< \mathbf{w}, \mathbf{x} \right>}{2} : ||\mathbf{w}|| \leq 1 
ight\}$ 

#### **Theorem**

 $\forall \gamma^* \in (0,1). \ \textit{H}_{\textit{lin}} \ \textit{is relaxed PACF learnable with} \\ g(\alpha,\gamma) = (\alpha \cdot \gamma - \gamma^*, \gamma^*) \ \textit{and sample and time complexities of} \\ poly(\frac{1}{\epsilon_{\gamma}}, \frac{1}{\epsilon_{\alpha}}, \frac{1}{\epsilon}, \log \frac{1}{\delta})$ 

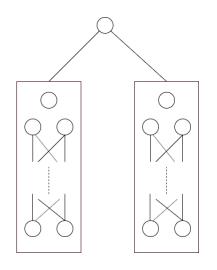
2. In logistic settings,

$$H_{\phi,L} \stackrel{\text{def}}{=} \{ \mathbf{x} \to \phi_{\ell}(\langle w, \mathbf{x} \rangle) : \mathbf{w} \in \mathcal{X}, \ell \in [0, L] \}$$

#### **Theorem**

 $\forall L>0, \forall \gamma^*\in (0,1), H_{lin}$  is relaxed PACF learnable with  $g(\alpha,\gamma)=(\alpha\cdot\gamma-\gamma^*,\gamma^*)$ , with sample and time complexities that are polynomial in  $poly(\frac{1}{\epsilon_\gamma},\frac{1}{\epsilon_\alpha},\frac{1}{\epsilon},\log\frac{1}{\delta})$ 

# Extension: 1.Embedding Individual Fairness Loss into General ML Training Objective



#### Formulation:

Given data set  $\mathcal{D} = (\mathbf{X}, \mathbf{y})$ ,  $\forall (\mathbf{x}_i, \mathbf{x}_j) \in M(\mathcal{D})$ ,  $h_1(\mathbf{x}_i) = \hat{y_i}, h_2(\mathbf{x}_j) = \hat{y_j}$ 

- 1. Utility Loss:  $\mathcal{L}_i(h_1(\mathbf{x}_i), y_i), \mathcal{L}_j(h_2(\mathbf{x}_j), y_j)$
- 2. Individual Fairness Loss:  $[d(h_1(\mathbf{x}_i), h_2(\mathbf{x}_j)) \Delta(\mathbf{x}_i, \mathbf{x}_j)]$  $\mathbb{1}(d > \Delta)$

For every possible pair, we back-propagate for both accuracy and individual fairness loss.

# 2. How to Get the Metric?

Right now, we are facing two problems:

- 1. How do we acquire the individual fairness metric  $\Delta$  for measuring the input distances?
- 2. If we have some subjective individual fairness notions available, can we elicit it into this model?
- Yes, both can be solved using metric learning. Suppose we are given some subjective notions of individual fairness notions, i.e.  $\forall x_i \in \mathcal{D}$ , there is an according label that is marked manually by experts (different from  $y_i$ ), then we can feed the instances and their labels into the general metric learning like large margin nearest neighbors<sup>1</sup>.

#### Question:

How is this architecture different from the Siamese Network?



<sup>&</sup>lt;sup>1</sup>Weinberger, K. Q., & Saul, L. K. (2009).

# **Algorithm 1:** Training process for the individually fair and accurate learning

Feed the training instances  $\mathcal X$  and the according subjective labels for individual fairness into the LMNN learning, get the distance metric  $\Delta$ ;

for 
$$1 \le i \le |\mathcal{X}|$$
 do  $|\mathcal{X}' = \mathcal{X}[i:] + \mathcal{X}[:i]$ 

1. for  $1 \le epoch \le 10$  do

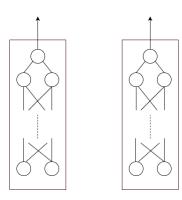
Feed  $\mathcal{X}, \mathcal{X}'$  into the architecture proposed. Calculate the prediction losses for sub-networks  $\mathcal{A}$  and  $\mathcal{B}$  separately and perform back-propagation.

#### end

- 2. Feed  $\mathcal{X}, \mathcal{X}'$  into the architecture proposed again.
- 3. Run through A and B, calculate the similarity loss and perform back-propagation.

#### end

#### 3. Alternative Structure



In the case where the fairness notions are hard binary labels: Just use two separate models. When optimizing the weights relative to loss, swap the outputs of the two individuals that are supposed to be treated similarly and then do back-propagation. This is suitable for the case that the subjective notions are binary.

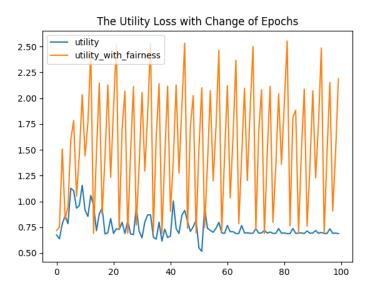
# 4.Experiment

- Data Set: Compas Recidivism Data <sup>2</sup>compas-scores-two-years. I use the binary values on whether the
  individuals commit a crime or not in the next few years as the
  labels. The rest of the features are inputs.
- Simulation of the Notions of the Subjective Individual
   Fairness: I generate a series of 0s and 1s for the individuals. If
   a pair of individual have the same notion, they are supposed to
   be treated similarly in terms of the machine prediction task.
- I define the individual fairness loss in the following way:

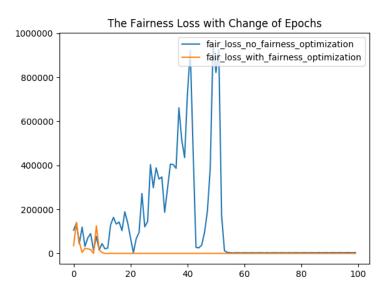
$$\mathcal{L}^F = \sum_{i=1}^{m-1} \sum_{j=i+1}^m [(\hat{y}_i - \hat{y}_j)^2] \mathbb{1}(i \text{ and } j \text{ should be treated similarly})$$

 $<sup>^2</sup> https://www.propublica.org/datastore/dataset/compas-recidivism-risk-score-data-and-analysis \\ + \Box > + \Box$ 

# 5. Experimental Results

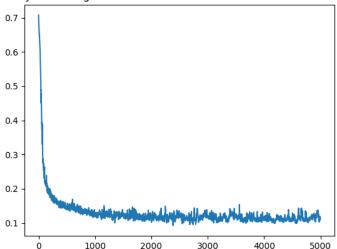


# 6.Experimental Results



# 7. Experimental Results

Utility loss change with no fairness constraint for 5000 combinations



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