

# Approximating likelihood ratios with calibrated classifiers

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Joint work with



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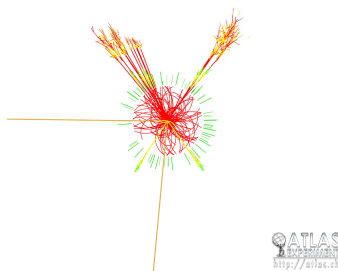


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See ([Cranmer et al., 2015](#)) for full details.

# Likelihood-free setup

- Complex simulator  $p$  parameterized by  $\theta$ ;
- Samples  $\mathbf{x} \sim p$  can be generated on-demand;
- ... but the likelihood  $p(\mathbf{x}|\theta)$  cannot be evaluated!



# Simple hypothesis testing

- Assume some observed data  $\mathcal{D} = \{\mathbf{x}_1, \dots, \mathbf{x}_n\}$ ;
- Test a null  $\theta = \theta_0$  against an alternative  $\theta = \theta_1$ ;
- The Neyman-Pearson lemma states that the most powerful test statistic is

$$\lambda(\mathcal{D}; \theta_0, \theta_1) = \prod_{\mathbf{x} \in \mathcal{D}} \frac{p_{\mathbf{x}}(\mathbf{x}|\theta_0)}{p_{\mathbf{x}}(\mathbf{x}|\theta_1)}.$$

- ... but neither  $p_{\mathbf{x}}(\mathbf{x}|\theta_0)$  nor  $p_{\mathbf{x}}(\mathbf{x}|\theta_1)$  can be evaluated!

# Straight approximation

1. Approximate  $p_{\mathbf{X}}(\mathbf{x}|\theta_0)$  and  $p_{\mathbf{X}}(\mathbf{x}|\theta_1)$  individually, using density estimation algorithms;
2. Evaluate their ratio  $r(\mathbf{x}; \theta_0, \theta_1)$ .

Works fine for low-dimensional data, but because of the curse of dimensionality, this is in general too difficult a problem!

$$\frac{p_{\mathbf{X}}(\mathbf{x}|\theta_0)}{p_{\mathbf{X}}(\mathbf{x}|\theta_1)} = r(\mathbf{x}; \theta_0, \theta_1)$$

*When solving a problem of interest, do not solve a more general problem as an intermediate step. – Vladimir Vapnik*

# Approximating likelihood ratios with classifiers

- Theorem: The likelihood ratio is invariant under the change of variable  $\mathbf{U} = s(\mathbf{X})$ , provided  $s(\mathbf{x})$  is monotonic with  $r(\mathbf{x})$ .

$$r(\mathbf{x}; \theta_0, \theta_1) = \frac{p_{\mathbf{X}}(\mathbf{x}|\theta_0)}{p_{\mathbf{X}}(\mathbf{x}|\theta_1)} = \frac{p_{\mathbf{U}}(s(\mathbf{x})|\theta_0)}{p_{\mathbf{U}}(s(\mathbf{x})|\theta_1)}$$

- Well, a classifier trained to distinguish  $\mathbf{x} \sim p_0$  from  $\mathbf{x} \sim p_1$  yields a decision function  $s(\mathbf{x})$  which is monotonic with  $r(\mathbf{x})$ .
- Estimating  $p(s(\mathbf{x})|\theta)$  is now easy, since the change of variable  $s(\mathbf{x})$  projects  $\mathbf{x}$  in a 1D space, where only the informative content of the ratio is preserved.
- Disentangle training from calibration.

# Inference and composite hypothesis testing

Approximated likelihood ratios can be used for inference, since

$$\begin{aligned}\hat{\theta} &= \arg \max_{\theta} p(\mathcal{D}|\theta) \\ &= \arg \max_{\theta} \prod_{\mathbf{x} \in \mathcal{D}} \frac{p(\mathbf{x}|\theta)}{p(\mathbf{x}|\theta_1)} \\ &= \arg \max_{\theta} \prod_{\mathbf{x} \in \mathcal{D}} \frac{p(s(\mathbf{x}; \theta, \theta_1)|\theta)}{p(s(\mathbf{x}; \theta, \theta_1)|\theta_1)}\end{aligned}\tag{1}$$

where  $\theta_1$  is fixed and  $s(\mathbf{x}; \theta, \theta_1)$  is a family of classifiers parameterized by  $(\theta, \theta_1)$ .

Accordingly, generalized (or profile) likelihood ratio tests can be evaluated in the same way.

# Parameterized learning

For inference, we need to build a family  $s(\mathbf{x}; \theta, \theta_1)$  of classifiers.

- One could build a classifier  $s$  independently for all  $\theta, \theta_1$ . But this is computationally expensive and would not guarantee a smooth evolution of  $s(\mathbf{x}; \theta, \theta_1)$  as  $\theta$  varies.
- Solution: build a single parameterized classifier instead, where parameters are additional input features (Baldi et al., 2016).

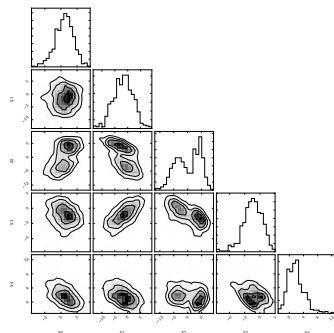
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 $\mathcal{T} := \{\};$   
while  $\text{size}(\mathcal{T}) < N$  do  
  Draw  $\theta_0 \sim \pi_{\Theta_0}$ ;  
  Draw  $\mathbf{x} \sim p(\mathbf{x}|\theta_0)$ ;  
   $\mathcal{T} := \mathcal{T} \cup \{((\mathbf{x}, \theta_0, \theta_1), y = 0)\}$ ;  
  Draw  $\theta_1 \sim \pi_{\Theta_1}$ ;  
  Draw  $\mathbf{x} \sim p(\mathbf{x}|\theta_1)$ ;  
   $\mathcal{T} := \mathcal{T} \cup \{((\mathbf{x}, \theta_0, \theta_1), y = 1)\}$ ;  
end while  
Learn a single classifier  $s(\mathbf{x}; \theta_0, \theta_1)$  from  $\mathcal{T}$ .
```



# Example: Inference from multidimensional data

Let assume 5D data  $\mathbf{x}$  generated from the following process  $p_0$ :

1.  $\mathbf{z} := (z_0, z_1, z_2, z_3, z_4)$ , such that  
 $z_0 \sim \mathcal{N}(\mu = \alpha, \sigma = 1)$ ,  
 $z_1 \sim \mathcal{N}(\mu = \beta, \sigma = 3)$ ,  
 $z_2 \sim \text{Mixture}(\frac{1}{2} \mathcal{N}(\mu = -2, \sigma = 1), \frac{1}{2} \mathcal{N}(\mu = 2, \sigma = 0.5))$ ,  
 $z_3 \sim \text{Exponential}(\lambda = 3)$ , and  
 $z_4 \sim \text{Exponential}(\lambda = 0.5)$ ;
2.  $\mathbf{x} := R\mathbf{z}$ , where  $R$  is a fixed semi-positive definite  $5 \times 5$  matrix defining a fixed projection of  $\mathbf{z}$  into the observed space.



Observed data  $\mathcal{D}$

Our goal is to infer the values  $\alpha$  and  $\beta$  based on  $\mathcal{D}$ .

🕒 Check out [\(Louppe et al., 2016\)](#) to reproduce this example.

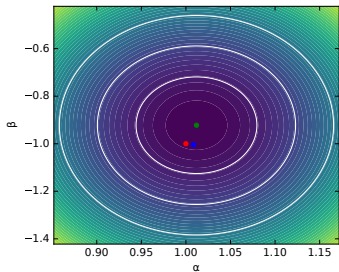
## Example: Inference from multidimensional data

Recipe:

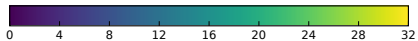
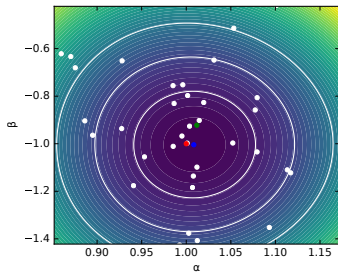
1. Build a single parameterized classifier  $s(\mathbf{x}; \theta_0, \theta_1)$ , in this case a 2-layer NN trained on 5+2 features, with the alternative fixed to  $\theta_1 = (\alpha = 0, \beta = 0)$ .
2. Find the approximated MLE  $\hat{\alpha}, \hat{\beta}$  by solving Eqn. 1.
  - Solve Eqn. 1 using likelihood scans or through optimization.
  - Since the generator is inexpensive,  $p(s(\mathbf{x}; \theta_0, \theta_1) | \theta)$  can be calibrated on-the-fly, for every candidate  $(\alpha, \beta)$ , e.g. using histograms.
3. Construct the log-likelihood ratio (LLR) statistic

$$-2 \log \Lambda(\alpha, \beta) = -2 \log \frac{p(\mathcal{D} | \alpha, \beta)}{p(\mathcal{D} | \hat{\alpha}, \hat{\beta})}$$

Exact  $-2 \log \Lambda(\alpha, \beta)$



Approx. LLR (smoothed by a Gaussian Process)



•  $\alpha = 1, \beta = -1$     • Exact MLE    • Approx. MLE

# Summary

- We proposed an approach for approximating LR in the likelihood-free setup.
- Evaluating likelihood ratios reduces to supervised learning. Both problems are deeply connected.
- Alternative to Approximate Bayesian Computation, without the need to define a prior over parameters.

# References

- Baldi, P., Cranmer, K., Faucett, T., Sadowski, P., and Whiteson, D. (2016). Parameterized Machine Learning for High-Energy Physics. *arXiv preprint arXiv:1601.07913*.
- Cranmer, K., Pavez, J., and Louppe, G. (2015). Approximating likelihood ratios with calibrated discriminative classifiers. *arXiv preprint arXiv:1506.02169*.
- Louppe, G., Cranmer, K., and Pavez, J. (2016). carl: a likelihood-free inference toolbox. <http://dx.doi.org/10.5281/zenodo.47798>, <https://github.com/diana-hep/carl>.

*Backup slides*

## Likelihood ratio of mixtures

# Diagnostics



# Applications in High Energy Physics