

# Estimating mutual information using sparse regression

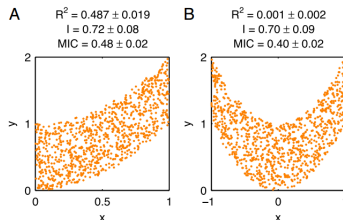
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(Joint work with Yuval Benjamini.)

# Mutual information (Shannon 1948)



- $I(X; Y) \in [0, \infty]$ . (0 if  $X \perp Y$ ,  $\infty$  if  $X = Y$  and  $X$  continuous.)
- Symmetry:  $I(X; Y) = I(Y; X)$ .
- Data-processing inequality

$$I(X; Y) \geq I(\phi(X); \psi(Y))$$

equality for  $\phi, \psi$  bijections

Image credit Kinney et al. 2014.

# Applications of $I(X; Y)$

- Feature selection (Peng et al. 2005, Fleuret 2004, Bennesar et al. 2015)
- Structure learning for graphical models using conditional mutual information  $I(X; Y|Z)$  (Vastano and Swinney 1988, Cheng et al. 1997, Bach and Jordan 2002)
- Quantifying information capacity of neurons

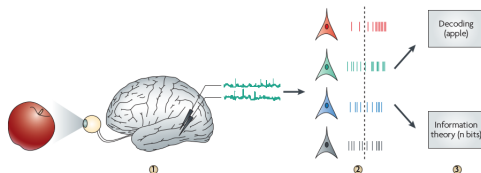


Image credits: Quiroga et al. (2009).

# How to estimate $I(X; Y)$

Suppose we observe pairs  $(X_i, Y_i)_{i=1}^n$  iid from density  $p(x, y)$

- Definition of mutual information:

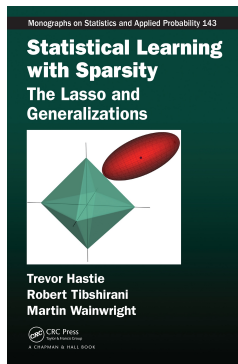
$$I(X; Y) = \int \log \left( \frac{p(x, y)}{p(x)p(y)} \right) p(x, y) dx dy$$

- Simply using plugging in kernel density estimate  $\hat{p}(x, y)$  leads to bias (Beirlant et al. 2001)
- Jackknifed estimate reduces bias (Ivanov and Rozhkova 1981)
- Alternative to KDE: nearest neighbors (Mnatsakanov et al. 2008, Goria et al. 2005, Singh et al. 2003)

# Problems in high dimensions

- Density estimation is known to have exponential complexity with respect to dimensionality.
- Many applications with high-dimensional  $X$ ,  $Y$ .
  - Gene expression time series
  - Functional magnetic resonance imaging
- One approach is to assume joint multivariate normality of  $X$ ,  $Y$ , but this reduces mutual information to a linear statistic.
- Other approaches: binning (Bialek et al. 1991, Paninski 2003), confusion matrix of a classifier (Treves 1997, Quiroga et al. 2009)

# Idea: Use sparsity!



- Suppose that  $Y \approx f(X) + \epsilon$ , where  $f$  depends *sparsely* on  $X$ .
- Can we exploit the sparsity to obtain an estimate of  $I(X; Y)$ ?

# Our proposal

Suppose we observe pairs  $(X_i, Y_i)_{i=1}^n$  iid from density  $p(x, y)$ .

- 1 Estimate a (sparse) regression model for  $\mathbf{E}[y|x]$ .
- 2 Estimate the *identification risk*  $p$  using cross-validation.
- 3 Use the identification risk to obtain a lower bound for the mutual information  $I(X; Y)$ :

$$I(X; Y) \geq g(p)$$

where  $g$  is a function that we derive theoretically.

# Multiple-response regression

- Pairs  $(x_i, y_i)_{i=1}^n$ , where  $X$  is  $p$ -dimensional and  $Y$  is  $q$ -dimensional.
- Data matrices  $\mathbf{X}_{n \times p}$ ,  $\mathbf{Y}_{n \times q}$ .
- For each column of  $Y$ , fit sparse model  $Y^{(i)} \approx X^T \beta^{(i)} + \epsilon$ , e.g. by using elastic net (Zou 2008),

$$\hat{\beta}^{(i)} = \operatorname{argmin}_{\beta} \|\mathbf{X}^T \beta^{(i)} - Y^{(i)}\|^2 + \lambda_2 \|\beta^{(i)}\|_2^2 + \lambda_1 \|\beta^{(i)}\|_1$$

- Or, fit a *random forest* model for each column of  $Y$  (Breiman 2001)



# Regression vs Identification loss

- Independent *test set*  $(x_i^*, y_i^*)_{i=1}^k$ .
- Use model to predict  $\hat{y}_i^* = (x_i^*)^T \hat{B}$  for  $i = 1, \dots, k$ .

Two ways to evaluate the predictive accuracy of the regression model:

- Regression (mean squared-error) loss:

$$\text{MSE} = \frac{1}{k} \sum_{i=1}^k \|y_i^* - \hat{y}_i^*\|^2.$$

- Identification loss (Kay 2008):

$$\text{IdLoss}_k = \frac{1}{k} \sum_{i=1}^k (1 - I\{\hat{y}_i^* \text{ is nearest neighbor of } y_i^*\}).$$

[note: point out that while idloss was introduced by Kay, that we are the first to consider theory, and add slide about 1d example/robustness]

# Identification loss and mutual information

- Define the identification risk as the expected identification loss

$$\text{IdRisk}_k = \mathbf{E}[\text{IdLoss}_k]$$

- Define the Bayes risk as the identification risk given the *true* model parameters. Hence,

$$\text{BayesRisk}_k \leq \text{IdRisk}_k.$$

- Theorem.** (Z., Benjamini 2016) There exists a function  $g_k$  such that

$$I(X; Y) \geq g_k(\text{BayesRisk}_k).$$

- Resulting estimator:

$$\hat{I}_{\text{IdLoss}}(X; Y) = g_k(\text{IdLoss}_k).$$

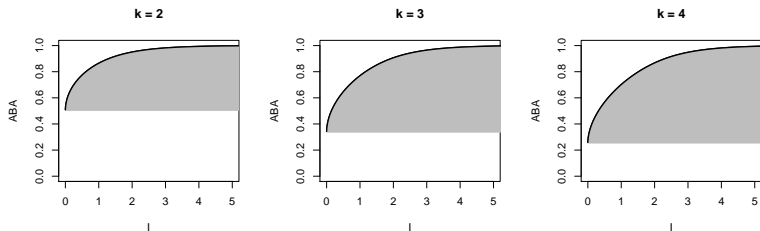
Leave- $k$ -out cross-validation (LkoCV) can be used for both squared-error loss and identification loss.

- Start with a dataset  $(x_i, y_i)_{i=1}^N$ .
- Let  $n = N - k$ . Consider all  $\binom{N}{k}$  partitions of the dataset into a test set  $(\mathbf{X}, \mathbf{Y})$  and training set  $(\mathbf{X}^*, \mathbf{Y}^*)$ .
- For each partition, compute the loss.
- Define the LkoCV loss as the average loss over  $\binom{N}{k}$  partitions.

*Computational note.* One can subsample to avoid computing all  $\binom{N}{k}$  partitions. In particular, if  $m = N/k$ , then one can use  $m$ -fold cross-validation which uses  $m$  partitions that have disjoint test sets.

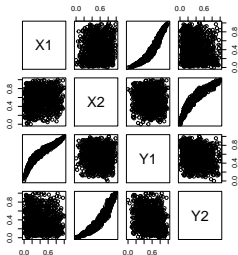
# Functions

Illustration of  $C_k = g_k^{-1}$



As information increases, the maximal identification risk goes to 0. [note: pictures need to be rotated]

# Simulation



- Generate data:  $(Y_1, Y_2) = f(X_1, X_2, \epsilon)$  where  $f$  is nonlinear.
- Add extra noise dimensions  $X_3, X_4, \dots$
- $n = 1000$ .
- Compare Nearest-Neighbor estimator (Mnatsakov et al, 2008, implemented in FNN) with our method using *Random Forest*.

# Simulation Results

True  $I(X; Y) = 4.615$ .

Extra dim	NN	RF $k = 10$	RF $k = 20$
0	<b>4.445</b>	3.989	3.924
1	3.040	<b>3.645</b>	3.610
2	1.773	<b>3.249</b>	3.182

## Section 2

# Theory

# Functional formulation

Bayes identification risk  $\text{BayesRisk}_k[p(x, y)]$  and mutual information  $I[p(x, y)]$  are both *functionals* of  $p(x, y)$ .

$$\text{BayesAcc}_k[p(x, y)] = \frac{1}{k} \int p_X(x_1) \dots p_X(x_k) \max_{i=1}^k p(y|x_i) dx_1 \dots dx_k dy.$$

$$I[p(x, y)] = \int p(x, y) \log \frac{p(x, y)}{p(x)p(y)} dx dy.$$

where  $\text{BayesAcc}_k = 1 - \text{BayesRisk}_k$ .



# Problem formulation

Take  $\iota > 0$ , and fix  $k \in \{2, 3, \dots\}$ . Let  $p(x, y)$  be a joint density (where  $(X, Y)$  could be random vectors of any dimensionality.) Supposing

$$I[p(x, y)] \leq \iota,$$

then can we find an upper bound,  $g_k^{-1}(\iota)$ , on  $\text{BayesAcc}_k[p(x, y)]$ ?

# Proof outline

- 1 Reduce problem to optimization over univariate densities.
- 2 Define the Lagrangian functional

$$\mathcal{L}[q(x)] = -\text{BayesAcc}_k[q(x)] + \lambda \int_0^1 q(x) dx + \nu I[q(x)]$$

which maps the univariate density  $q(x)$  to a real number.

- 3 Compute the functional derivative of the Lagrangian

$$\nabla \mathcal{L}[q](x) = -t^{k-1} + \lambda + \nu(1 + \log q(x))$$

- 4 Set  $\nabla \mathcal{L}[q](x) = 0$ , yielding

$$q^*(t) = \alpha e^{\beta t^{k-1}}.$$

- 5 Check that local minimizer is global minimizer.

**Theorem.** For any  $\iota > 0$ , there exists  $\beta_\iota \geq 0$  such that defining

$$q_\beta(t) = \frac{\exp[\beta t^{k-1}]}{\int_0^1 \exp[\beta t^{k-1}]},$$

we have

$$\int_0^1 q_{\beta_\iota}(t) \log q_{\beta_\iota}(t) dt = \iota.$$

Then,

$$\sup_{I(X;Y)=\iota} \text{BayesAcc}_k = \int_0^1 q_{\beta_\iota}(t) t^{k-1} dt = g_k^{-1}(\iota).$$

## Section 3

# Conclusion

# Application to gene expression time series

to be contd

# Related work and future directions

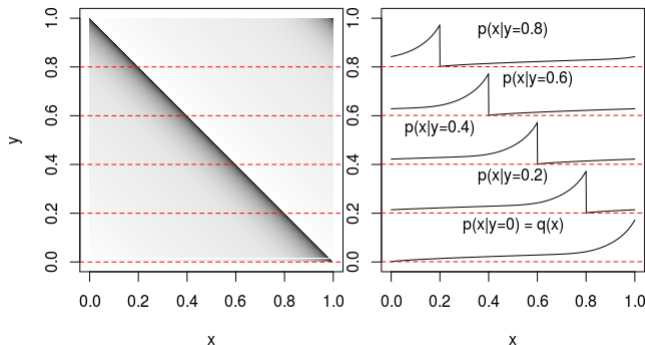
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## Section 4

The End

# Reduced Problem

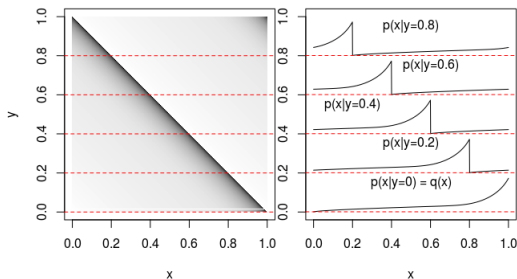
Rather than show the whole proof, we consider a simplified problem to illustrate the methods.



Actually, the simplified problem is equivalent to the full problem and we get the same answer (but this is non-trivial).



# Reduced Problem



- $p(x, y)$  on unit square with uniform marginals.
- The conditional distributions  $p(x|y)$  are just “shifted” copies of a common density,  $q(x)$ , on  $[0, 1]$

$$p(x|y) = q(x - y + I\{x < y\})$$

- Furthermore,  $q(x)$  is increasing in  $x$ .

The information and average Bayes error can be written in terms of  $q(x)$ .

$$I[p(x, y)] = \int_0^1 q(x) \log q(x) dx$$

$$\text{BayesAcc}_k[p(x, y)] = \int_{[0,1]^k} \max_{i=1}^k q(x_i) dx_1 \cdots dx_k$$

Overload the notation and “redefine” information and average Bayes error as functionals of  $q(x)$ .

$$I[q(x)] \stackrel{\text{def}}{=} \int_0^1 q(x) \log q(x) dx$$

$$\text{BayesAcc}_k[q(x)] \stackrel{\text{def}}{=} \frac{1}{k} \int_{[0,1]^k} \max_{i=1}^k q(x_i) dx_1 \cdots dx_k$$

# Optimization problem

We now pose the question: how do we find  $q(x)$  which maximizes  $\text{BayesAcc}_k[q(x)]$  subject to  $I[q(x)] \leq \iota$ ?

- *Domain of the optimization:* Recall that  $q(x)$  satisfies  $q(x) \geq 0$ ,  $\int_0^1 q(x)dx = 1$ , and is increasing in  $x$ . Let  $\mathcal{Q}$  denote the space of functions on  $[0, 1] \rightarrow [0, \infty)$  which are increasing in  $x$ .
- *Constraints:* We have two remaining constraints,  $I[q(x)] \leq \iota$  and  $\int_0^1 q(x)dx = 1$ .

Hence the problem is

$$\text{maximize}_{q(x) \in \mathcal{Q}} \text{BayesAcc}_k[q(x)] \text{ subject to } \int_0^1 q(x)dx = 1 \text{ and } I[q(x)] \leq \iota.$$

# Optimization problem

maximize $_{q(x) \in \mathcal{Q}}$  BayesAcc $_k[q(x)]$  subject to  $\int_0^1 q(x)dx = 1$  and  $I[q(x)] \leq \iota$ .

- Does a solution exist? Yes, because the space of measures with density  $q(x)$  satisfying  $I[q(x)] \leq \iota$  is tight, and both the constraints and objective are continuous wrt to the topology of weak convergence.
- Given a solution  $q^*(x)$  exists, there exist Lagrange multipliers  $\lambda \in \mathbb{R}$  and  $\nu > 0$  such that  $q^*$  minimizes

$$\begin{aligned}\mathcal{L}[q(x)] &= -\text{BayesAcc}_k[q(x)] + \lambda \int_0^1 q(x)dx + \nu I[q(x)] \\ &= \int_0^1 (-t^{k-1} + \lambda + \nu \log q(x))q(x)dx.\end{aligned}$$

# Functional derivatives

- Taylor expansions are a useful trick for computing functional derivatives
- We can compute the functional derivative of  $\mathcal{L}[q(x)]$  by writing

$$\begin{aligned}\mathcal{L}[q(x) + \epsilon \xi(x)] &= \int_0^1 (-t^{k-1} + \lambda + \nu \log(q(x) + \epsilon \xi(x)))(q(x) + \epsilon \xi(x)) dx. \\ &\approx \int (q(x) + \epsilon \xi(x))(-t^{k-1} + \lambda + \nu \{\log q(x) + \frac{\epsilon \xi(x)}{q(x)}\}) dx \\ &\approx \mathcal{L}[q(x)] + \int_0^1 (-t^{k-1} + \lambda + \nu(1 + \log q(x))) \epsilon \xi(x) dx.\end{aligned}$$

- Hence

$$\nabla \mathcal{L}[q](x) = -t^{k-1} + \lambda + \nu(1 + \log q(x))$$

# Variational magic!

Suppose we set the functional derivative to 0,

$$0 = \nabla \mathcal{L}[q](t) = -t^{k-1} + \lambda + \nu + \nu \log q(t).$$

Then we conclude that the optimal  $q^*(t)$  takes the form

$$q^*(t) = \alpha e^{\beta t^{k-1}}$$

for some  $\alpha > 0$ ,  $\beta > 0$ .

From the constraint  $\int q(t) dt = 1$ , we get

$$q_\beta(t) = \frac{e^{\beta t^{k-1}}}{\int e^{\beta t^{k-1}} dt}.$$

**Theorem.** For any  $\iota > 0$ , there exists  $\beta_\iota \geq 0$  such that defining

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we have

$$\int_0^1 q_{\beta_\iota}(t) \log q_{\beta_\iota}(t) dt = \iota.$$

Then,

$$\sup_{I(X;Y)=\iota} \text{BayesAcc}_k = \int_0^1 q_{\beta_\iota}(t) t^{k-1} dt = g_k^{-1}(\iota).$$