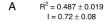
Estimating mutual information for high-dimensional sparse relationships

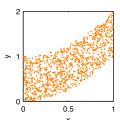
Charles Zheng

Stanford University

January 14, 2017

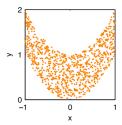
(Joint work with Yuval Benjamini.)





B
$$R^2 = 0.001 \pm 0.002$$

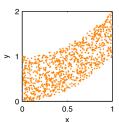
I = 0.70 ± 0.09



Mutual information $I(\vec{X}; \vec{Y})$

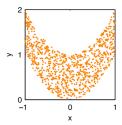
• measures dependence between two random vectors, \vec{X} and \vec{Y}

A $R^2 = 0.487 \pm 0.019$ I = 0.72 ± 0.08



B
$$R^2 = 0.001 \pm 0.002$$

I = 0.70 ± 0.09

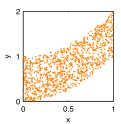


Mutual information $I(\vec{X}; \vec{Y})$

- measures dependence between two random vectors, \vec{X} and \vec{Y}
- applies to nonlinear and multidimensional relationships (unlike correlation)

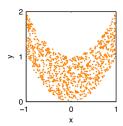
A
$$R^2 = 0.487 \pm 0.019$$

I = 0.72 ± 0.08



B
$$R^2 = 0.001 \pm 0.002$$

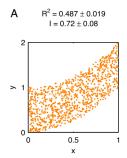
I = 0.70 ± 0.09



Mutual information $I(\vec{X}; \vec{Y})$

- measures dependence between two random vectors, \vec{X} and \vec{Y}
- applies to nonlinear and multidimensional relationships (unlike correlation)
- is difficult to estimate in high dimensions

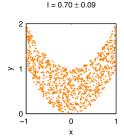
В



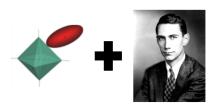
Mutual information $I(\vec{X}; \vec{Y})$

- measures dependence between two random vectors, \vec{X} and \vec{Y}
- applies to nonlinear and multidimensional relationships (unlike correlation)
- is difficult to estimate in high dimensions

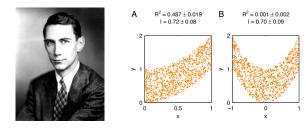
We combine machine learning (sparse estimation) with information theory to obtain better estimates of $I(\vec{X}; \vec{Y})$



 $B^2 = 0.001 + 0.002$



Mutual information I(X; Y)



Introduced in Shannon's 1948 paper, "A mathematical theory of communication"

Image credit Kinney et al. 2014.

Applications of I(X; Y)

Mutual information has since been applied to many areas outside of information theory

Applications of I(X; Y)

Mutual information has since been applied to many areas outside of information theory

Applications [edit]

In many applications, one wants to maximize mutual information (thus

- In search engine technology, mutual information between phrases
- . In telecommunications, the channel capacity is equal to the mutua
- Discriminative training procedures for hidden Markov models have
- RNA secondary structure prediction from a multiple sequence alig
- Phylogenetic profiling prediction from pairwise present and disapp
- Mutual information has been used as a criterion for feature selectithe minimum redundancy feature selection.
- . Mutual information is used in determining the similarity of two diffe
- Mutual information of words is often used as a significance functio words; rather, one counts instances where 2 words occur adjacen another, goes up with N.
- Mutual information is used in medical imaging for image registratic reference image, this image is deformed until the mutual information
- · Detection of phase synchronization in time series analysis
- . In the infomax method for neural-net and other machine learning,

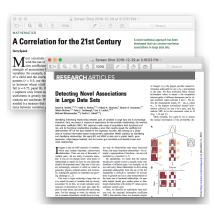
Engineering, biology, computer science, physics, medicine

Comparing I(X; Y) with Pearson correlation



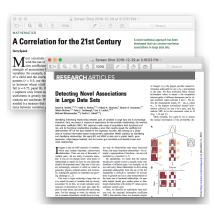
 In many applications scientists are interested in dependence, not correlation (Reshef et al. 2011, Speed 2011).

Comparing I(X; Y) with Pearson correlation



- In many applications scientists are interested in dependence, not correlation (Reshef et al. 2011, Speed 2011).
- Only mutual information (and derived quantities) measures dependence directly

Comparing I(X; Y) with Pearson correlation



- In many applications scientists are interested in dependence, not correlation (Reshef et al. 2011, Speed 2011).
- Only mutual information (and derived quantities) measures dependence directly

Problems with mutual information

- Hard to interpret (compared to R^2)
- Hard to estimate (compared to R^2)

Can we make I(X; Y) easier to interpret?

• Define the "informational correlation" (Linfoot 1957)

$$Cor_{Info}(X, Y) = \sqrt{1 - e^{-2I(X;Y)}}$$

Can we make I(X; Y) easier to interpret?

• Define the "informational correlation" (Linfoot 1957)

$$Cor_{Info}(X,Y) = \sqrt{1 - e^{-2I(X;Y)}}$$

- Then $Cor_{Info}(X, Y) \in [0, 1]$.
- For (X, Y) bivariate normal,

$$|\mathsf{Cor}_{\mathsf{Pearson}}(X,Y)| = \mathsf{Cor}_{\mathsf{Shannon}}(X,Y)$$

Can we make I(X; Y) easier to interpret?

• Define the "informational correlation" (Linfoot 1957)

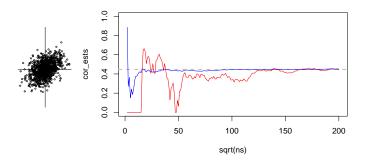
$$Cor_{Info}(X,Y) = \sqrt{1 - e^{-2I(X;Y)}}$$

- Then $Cor_{Info}(X, Y) \in [0, 1]$.
- For (X, Y) bivariate normal,

$$|\mathsf{Cor}_{\mathsf{Pearson}}(X,Y)| = \mathsf{Cor}_{\mathsf{Shannon}}(X,Y)$$

Difficulty of estimating I(X; Y)

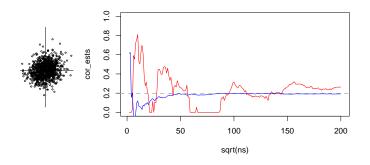
Example with $Cor_{Pearson}(X, Y) = Cor_{Info}(X, Y) = 0.44$.



8 / 49

Difficulty of estimating I(X; Y)

Example with $Cor_{Pearson}(X, Y) = Cor_{Info}(X, Y) = 0.2$.



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:

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

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:

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

- Kernel density estimate approaches estimate p(x, y) (Beirlant et al. 2001, Ivanov and Rozhkova 1981)
- Nearest neighbor estimators rely on distance-based computations (Mnatsakanov et al. 2008, Goria et al. 2005, Singh et. al. 2003)

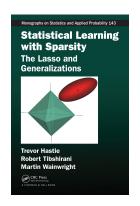
- Density estimation is known to have *exponential complexity* with respect to dimensionality.
 - E.g. to get the same precision, you need 10 observations for univariate X, Y but 1000 for trivariate \vec{X}, \vec{Y} .

- Density estimation is known to have exponential complexity with respect to dimensionality.
 - E.g. to get the same precision, you need 10 observations for univariate X, Y but 1000 for trivariate \vec{X}, \vec{Y} .
- Many applications with high-dimensional X, Y.
 - Gene expression time series
 - Functional magnetic resonance imaging

- Density estimation is known to have *exponential complexity* with respect to dimensionality.
 - E.g. to get the same precision, you need 10 observations for univariate X, Y but 1000 for trivariate \vec{X}, \vec{Y} .
- 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.

- Density estimation is known to have exponential complexity with respect to dimensionality.
 - E.g. to get the same precision, you need 10 observations for univariate X, Y but 1000 for trivariate \vec{X}, \vec{Y} .
- 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)

First 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)?

Second idea: link prediction accuracy to mutual information

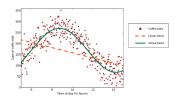
• If I(X; Y) > 0, then X carries information about Y and vice-versa.

Second idea: link prediction accuracy to mutual information

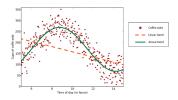
- If I(X; Y) > 0, then X carries information about Y and vice-versa.
- Therefore, we can predict Y from X (or X from Y)

Second idea: link prediction accuracy to mutual information

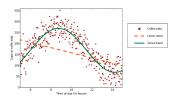
- If I(X; Y) > 0, then X carries information about Y and vice-versa.
- Therefore, we can predict Y from X (or X from Y)
- We know that often prediction accuracy implies a lower bound for mutual information (e.g. Fano 1952)



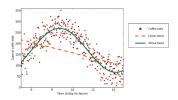
• Suppose you observe $(\vec{X}^{(i)}, Y^{(i)})_{i=1}^n$ where $Y^{(i)} = f(\vec{X}^{(i)}) + \epsilon$, where f is an unknown function and ϵ is noise. (Also, assume $\mathbf{E}[\epsilon] = 0$.)



- Suppose you observe $(\vec{X}^{(i)}, Y^{(i)})_{i=1}^n$ where $Y^{(i)} = f(\vec{X}^{(i)}) + \epsilon$, where f is an unknown function and ϵ is noise. (Also, assume $\mathbf{E}[\epsilon] = 0$.)
- ullet The goal in regression is to recover the unknown function f.



- Suppose you observe $(\vec{X}^{(i)}, Y^{(i)})_{i=1}^n$ where $Y^{(i)} = f(\vec{X}^{(i)}) + \epsilon$, where f is an unknown function and ϵ is noise. (Also, assume $\mathbf{E}[\epsilon] = 0$.)
- The goal in regression is to recover the unknown function f.
- In *linear regression*, we assume *f* is linear.



- Suppose you observe $(\vec{X}^{(i)}, Y^{(i)})_{i=1}^n$ where $Y^{(i)} = f(\vec{X}^{(i)}) + \epsilon$, where f is an unknown function and ϵ is noise. (Also, assume $\mathbf{E}[\epsilon] = 0$.)
- The goal in regression is to recover the unknown function f.
- In *linear regression*, we assume *f* is linear.
- if we do not assume a particular form for f, we can use nonparametric regression.

Background: Sparse regression

• When \vec{X} is high dimensional, classical regression techniques perform poorly.

Background: Sparse regression

- When \vec{X} is high dimensional, classical regression techniques perform poorly.
- If the true function f only depends on a small number of components in \vec{X} , we can still do well if we use *sparse* regression methods.

Background: Sparse regression

- When \vec{X} is high dimensional, classical regression techniques perform poorly.
- If the true function f only depends on a small number of components in \vec{X} , we can still do well if we use *sparse* regression methods.

	Classical	Sparse
Linear	Ordinary Least-Squares	Elastic net
	(Gauss 1975?)	(Zou 2008)
Nonpar.	LOWESS	Random forests
	(Cleveland 1979)	(Breiman 2001)

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 $\mathbf{E}[y|x]$.
- Assess the prediction accuracy of the model using identification risk
- ① Use the identification risk to obtain a lower bound for the mutual information I(X; Y)

Multiple-response regression

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

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

• 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, ..., k.

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

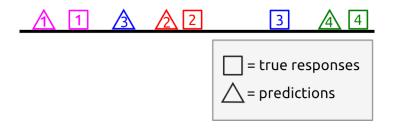
• Regression (mean squared-error) loss:

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

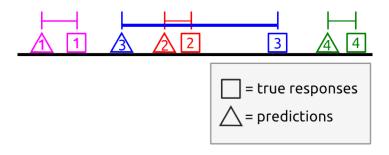
Identification loss (Kay 2008):

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

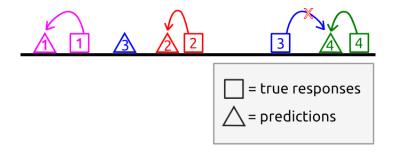
Regression vs Identification loss



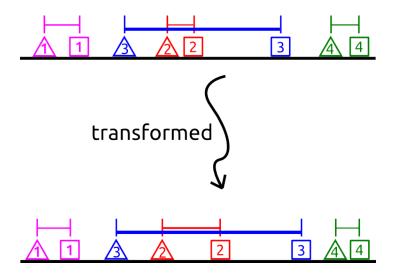
Mean-squared error



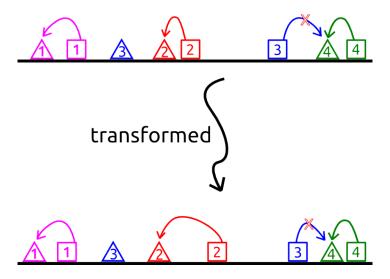
Identification loss



Mean-squared error changes under nonlinear scaling



Identification loss robust under nonlinear scaling



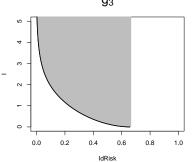
Identification loss and mutual information

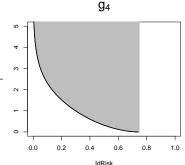
Define the identification risk as the expected identification loss

$$IdRisk_k = \mathbf{E}[IdLoss_k]$$

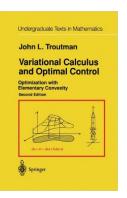
• **Theorem.** (Z., Benjamini 2017) There exists a function g_k such that

$$I(X;Y) \geq g_k(\mathsf{IdRisk}_k).$$
g $_3$



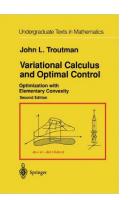


We use calculus of variations to obtain this result.



We use calculus of variations to obtain this result.

• Mutual information is a functional of p(x, y).



$$I[p(x,y)] = \mathbf{E}\left[\log \frac{p(X,Y)}{p(X)p(Y)}\right].$$

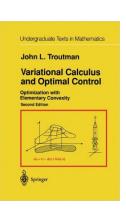
We use calculus of variations to obtain this result.

• Mutual information is a functional of p(x, y).

$$I[p(x,y)] = \mathbf{E}\left[\log \frac{p(X,Y)}{p(X)p(Y)}\right].$$

 Identification risk is lower-bounded by another functional—the Bayes risk.

$$\mathsf{BayesRisk}_k[p(x,y)] = 1 - \mathbf{E}[\max_{i=1}^k p(Y|X_i)].$$



We use calculus of variations to obtain this result.

• Mutual information is a functional of p(x, y).

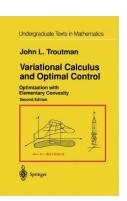
$$I[p(x,y)] = \mathbf{E}\left[\log \frac{p(X,Y)}{p(X)p(Y)}\right].$$

 Identification risk is lower-bounded by another functional—the Bayes risk.

$$\mathsf{BayesRisk}_k[p(x,y)] = 1 - \mathbf{E}[\max_{i=1}^k p(Y|X_i)].$$

• $g_k(u) = \inf_{p(x,y)} I[p(x,y)]$

subject to BayesRisk $_k[p(x, y)] \ge u$.



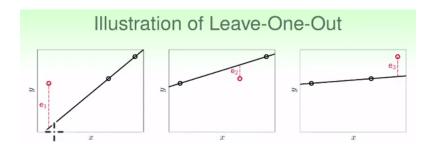
Our proposal

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

- **1** Estimate a (sparse) regression model for $\mathbf{E}[y|x]$.
- Compute identification loss, IdLossk, using leave-k-out.
- Stimate mutual information using

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

What is leave-k-out cross-validation?



- Randomly hold out a subset of size k.
- Use remaining data to predict the held-out data.
- Obtain the average prediction error.

Image credit Hsuan-Tien Lin

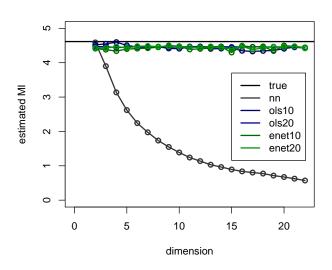
Section 2

Applications

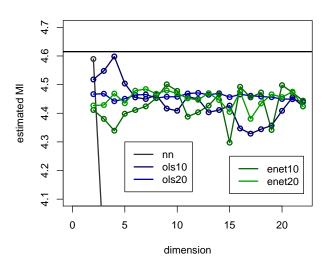
Simulation

- Generate data: $(Y_1, Y_2) = (X_1, X_2)^T B + \epsilon$ where B is a randomly generated coefficient matrix.
- Add extra noise dimensions X_3, X_4, \ldots
- n = 1000.
- Compare Nearest-Neighbor estimator (Mnatsakov et al, 2008, implemented in FNN) with our method using OLS and elastic net (sparse).

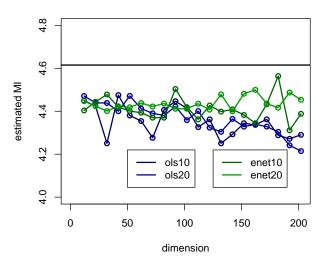
Simulation Results - I. low dimension



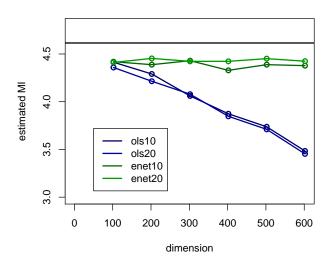
Simulation Results - I. low dimension



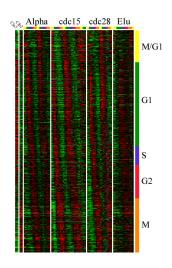
Simulation Results - II. medium dimension



Simulation Results - III. high dimension



Application to gene expression time series



- Data from Spellman et al. 1998
- Expression levels of 6178 yeast genes during cell cycle
- Total 73 time points per gene

Groups of genes

Group	No. genes
unknown	396
cell cycle	27
DNA replication	27
transport	19
cytoskeleton	17
chromatin structure	16

Total 145 different categories (only top 6 shown).

Canonical correlations between time series

Top canonical correlation (Hotelling 1936)

	CC	DR	Tr	Су	CS
CC		1	1	1	1
DR			1	0.99	0.99
Tr				0.99	0.98
Су					0.98
CS					

 $CC = cell\ cycle,\ DR = DNA\ replication,\ Tr = transport,$ $Cy = cytoskeleton,\ CS = chromatin\ structure$

Sparse canonical correlations between time series

Using sparse CCA* (Witten and Tibshirani 2009).

	CC	DR	Tr	Су	CS
CC		0.93	0.79	0.88	0.87
DR			0.76	0.76	0.91
Tr				0.65	0.68
Су					0.83
CS					

$$CC = cell\ cycle,\ DR = DNA\ replication,\ Tr = transport,$$
 $Cy = cytoskeleton,\ CS = chromatin\ structure$

^{*:} Default settings in R package PMA

Information correlations between time series

Taking the max of $\hat{I}(X; Y)$ and $\hat{I}(Y; X)$.

	CC	DR	Tr	Су	CS
CC		0.92	0.79	0.97	0.81
DR			0.85	0.90	0.92
Tr				0.73	0.71
Су					0.93
CS					

 $CC = cell\ cycle,\ DR = DNA\ replication,\ Tr = transport,$ $Cy = cytoskeleton,\ CS = chromatin\ structure$

Related work and future directions

to be contd

Section 3

The End

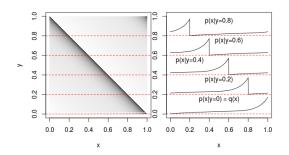
Reduced Problem

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.

Simplified formulae

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

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

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

Simplified formulae

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

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

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

Optimization problem

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

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

Hence the problem is

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

Optimization problem

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

- Does a solution exist? Yes, because the space of measures with density q(x) satisfying $I[q(x)] \le \iota$ 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

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

Functional derivatives

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

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

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

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

Result

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

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

we have

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

Then,

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