## NOTES DURING PHD

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#### 1. Meetings.

1.1. Nonnegative martingales and E-process.

25.08.2025. Practical goal:

- Compare speed of conditional E in BiasedUrn and R: about 3x performance, not worth it to import
- Add save E into the package: forked from

This week's goal theory-wise

- Just write down again the KL for Gaussian family for repeatition.
- Restate the connection between integral and sequential product of UI: haven't found the literature yet
- Try to re-state the simple and anti-simple case
- What is the seq-RIPr and seq-COND?
- Why do we need to have a general KL measure in general paper
- Read carefully how the maximum is found in the log-optimal paper
- · Reproduce with safestats packages for the
- UMP in math. stat. course lecture notes.
- I wanna write down  $d\nu$  and dX and all that.

The difference between the simple and anti-simple case In short, the simple case is where  $\Sigma_q - \Sigma_p$  is negative which mean we can find a RIPr via a prior (or a element) of P.

15.08.2025. Sebastian did a great presentation regarding testing quantile given filtrations. I would formulate here and also add a picture.

The question in mind is to test whether data X is from a hypothesis  $\mathcal H$  that is non-parametric:

$$\mathcal{H} = \{ P \in \mathcal{P}(X) \mid \mathbb{E}_p \left[ \phi_i(X) = 0 \right], \beta = 1, 2, \dots, d \}$$
 where  $\mathcal{P}(X)$  is all distribution on  $X$ .

The simplest instance where testing X with the same mean  $\mu$  where  $\phi_1(X) = X - \mu$ . Larson has proven that the 'optimal' E-variable must be in the form:

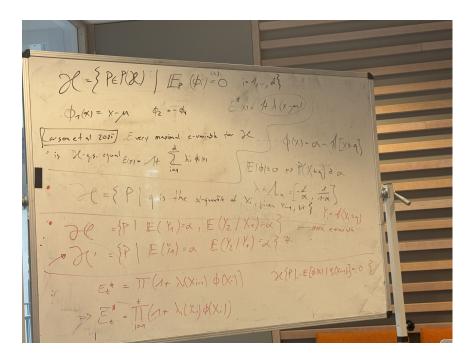
$$S := 1 + \sum_{i=1}^{d} \lambda_i \phi_i(X)$$

Sebastian is mainly interested in the cases if there is any gain in  $\mathbb{E}$ -variables compared to a coraser filtrations. Imagine we have conditioned on the original data  $X_i$ ,  $i=1,2,\ldots$ , you would think that the hypothesis case constructed as below:

$$(1.1) \mathcal{H} = \{ P \in \mathcal{P}(X) \mid \mathbb{E}[Y_2 \mid X_1] = \alpha \}$$

$$\mathcal{H}' = \{ P \in \mathcal{P}(X) \mid \mathbb{E}[Y_2 \mid Y_1] = \alpha \}$$

where  $Y_i = 1_{X_i \leq q}$ . He showed that both  $\mathcal{H}$  and  $\mathcal{H}'$  are convex and  $\mathcal{H}$ 's closed convex hull is not the same, otherwise it would be kind of pointless.



14.08.2025. This is a short discussion regarding UI, specifically about the difference between prequential and integral representation.

Supposed iid data  $\mathcal{D} = \{X_1, X_2, \dots, X_n\}$  and we denote the data up to time i by  $X^{(i)} = \{X_1, X_2, \dots, X_i\}$ .

One way to instantiate UI by

$$\frac{\prod_{i=1}^{n} P_{\tilde{\theta}_{alt|i-1}}(X_i)}{P_{\hat{\theta}_0}(X^{(n)})}$$

where  $\tilde{\theta}_{alt|i-1}$  is any estimator in alternative based on the first i-1 data points  $X^{(i-1)}$  and  $\hat{\theta}_0 = \arg\max_{\theta \in \Theta_0} \prod_i p_{\theta}(X_i)$  is the MLE estimator under the null. See more details in section 7 of UI paper.

Notice that the sequence of the data  $\mathcal{D} = \{X_1, \dots X_n\}$  really matters. Imagine you obtian  $\mathcal{D}'$  with a rearranged sequence and thus slightly different  $\tilde{\theta}$  and hence slightly different value at the denominator.

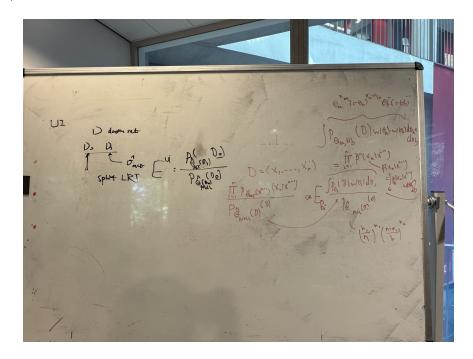
Revisit factorization of probability, it's also called chain rule or general product rule:

$$P(X_{1}, X_{2},..., X_{n}) = P(X_{1}) P(X_{2} | X_{1}) \cdots P(X_{n} | X_{1}, X_{2},..., X_{n-1})$$

$$= \prod_{i=1}^{n} P(X_{i} | X^{(i-1)})$$

QUESTION 1.1. Below is a mixture of all nonnegative test martingale/e-process?

$$\frac{\int p_{\theta}(X^{(n)})w(\theta)\,\mathrm{d}\theta}{P_{\hat{\theta}_{\text{null}}}(X^{(n)})}$$



25.07.2025. I only briefly went over the conformal prediction and fisher's noncentral hypergeometric distribution. There were some discussions on what the conformal prediction is.

Imagine you have a classifier for images (dogs, cat, etc.) and is trained via N datasets. For the next prediction, we need something to quantify the uncententy to say that

The label for  $X_{N+1}$  I gave being X (here could be any label), has

Peter gave a algorithmic explanantion where the predicted labels are gaven for each label, then run through against the previous training datasets. Rank them, cut off the tailing  $\alpha$  percent then we can say we are confident about our prediction with  $1-\alpha$ .

However, this is awfully similar to the permutation test, by Sebastian. Yeah it does look a lot like just ranking the prediction and give a p-value.

Alexander also suggested using Gaussian for conformal prediction might be too confusing as the parameter and the prediction kind of just are the same thing. Maybe try Poisson example where parameter is in real number while the prediction is in  $\mathbb{Z}$ .

What I do not follow is the output of the classifier is a weighted matrix over all the labels. What is the ranking being done over? Is it

**2. TODO.** This is just a playground for me! No need to be too nicely formatted.

#### 2.1. Questions.

- What is the regularity condition in universal inference condition?
- Why do we want to have data-dependent significant level  $\alpha$
- What is the difference between p-process and E-process and nonnegative martingale?
- What is a adapted sequence of random sets, random variables
- Wald's sequential likelihood ratio test
- What is Radon-Nikodym derivative
- Try to explain post-hod and ad-hoc in plain English
- What would happen if you just multiply E-variables together? Shouldn't you consider the
- What is the evidence in testing? Is likelihood ratio the best one we have for evidence

- What is the difference between carrying measure and carrying measure?
- Write down the lemma/fact that each element in a regular exponential family is continuous with respect to each other and could be used as carrier. Is it a one-to-one mapping?
- Mean, Variance, Fisher information
- Mean-value Parameterization Convexity of mean-value spaces and canonical spaces
- Duality
- · Loewner ordering
  - 2.2. Filtrations and sigma-fields.

DEFINITION 2.1 (Filtration). Let  $(\Omega, \mathcal{A}, P)$  be a probability space. An increasing sequence  $\mathcal{F}_n$  of sub- $\sigma$  algebras of  $\mathcal{A}$  (i.e.  $\mathcal{F}_0 \subseteq \mathcal{F}_1 \subseteq \cdots \subseteq \mathcal{A}$ ) is called a filtration.

The sub- $\sigma$ -algebra is just the sigma algebra of the  $X_0, X_1, X_n$ .

What is the difference between the stopping time and random times that can take possibly infinity.

Filtration sigma-field at time t, filtration is a n increasing sequence of sigma-fields.

What is the  $\limsup$  here?  $A_t$  is an adapted sequences of events in some filtered probability space.

$$A_{\infty} := \limsup_{t \to \infty} A_t := \bigcap_{t \in \mathbb{N}} \bigcup_{s > t} A_s$$

DEFINITION 2.2 (Absolute continuous). Let p,q be two probability distributions. p is called *absolutely continuous* with respect to q if the

More explicitly, this reads  $(p \ll q)$  that

$$q(A) = 0 \quad \Rightarrow \quad p(A) = 0, \quad \text{for any } A \in \mathcal{F}_t \text{ and } t \in \mathbb{N}.$$

2.3. *All the inequalities*. Hoeffding, Bernstein, McDiarmid, Talagrand's Evidence lower bound

line crossing inequality average treatment effect mixture adaptive design

- 2.4. Papers, talks, textbook, and more topics. Clarke Barron 1990 1994 about
- · Stein's Methods
- · Ito's process
- Levy's process
- Gaussian process
  - 2.5. Basic Concepts.

DEFINITION 2.3 (Admissibility). In E-process, a process is inadmissible if there exists a process that is strictly better than it at certain time points.

Calibration

**Dominates** 

UMP

Understanding the Fourier transform in terms of transform Placing a restriction on the Fourier transform == smaller function space Minimax setup

Integration / Measure

Topology crash course

The topologist Stephen Smale stunned the mathematical community in 1958 [Sma58] when he proved it was possible to turn the sphere inside out without introducing any creases. Several ways to do this are beautifully illustrated in video recordings [Max77, LMM94, SFL98].

Tower properties

**3. TLDR for papers.** Here I really hope that I can somehow summarize in my words of the talks or presentation I have been to or literature I have read into a *short* paragraph.

There are only two requirement:

- short: ideally in one sentence, but it should definitely shorter than the abstract
- memorable/impressionable: for example, instead of precisely stating the LLN, it is better to write "keep tossing a fair coin, we are sure about the average being 0.5".

Admissible Anytime-Valid Sequential Inference Must Rely on Nonnegative Martingale (Ramdas et al., 2022a):

p-process, e-process, stopping times are sort of the same: we can always find a nonnegative martingale for them. However, there is a gap.

The numeraire e-variable and reverse information projection (Larsson, Ramdas and Ruf, 2024):

Shows that there is always a numeraire  $\mathbb{E}$ -variable  $X^*$  in the form of  $\mathbb{E}_Q\left[X/X^*\right] \leq 1$  for every  $\mathbb{E}$ -variable X in the simple alternative Q vs composite null  $\mathcal{P}$  setting. It is unique (up to Q-nullset), log-optimal, and connection between the effective null!

Reverse Information Projections and Optimal E-statistics (Lardy, Grünwald and Harremoës, 2024):

There is always a RIPr (?) even when the KL is infinite in the simple alternative Q vs composite null  $\mathcal{P}$  setting. In simple words, if the KL is infinite (bad at describing compared), there is still a relatively better version to describe that!.

Is the only minus infty or zero ?x

The extended Ville's inequality for nonintegrable nonnegative supermartingales

#### 4. Reverse Inverse Projections (RIPr).

4.1. Li's Algorithm. Originally developed by Li (1999), the following algorithm is written based on the notes from Grünwald, de Heide and Koolen (2024a); Hao et al. (2024). To quote Brinda (2018), Li's inequality requires the family  $\mathcal Q$  to have a uniformly bounded density ratio.

Li obtains the RIPr in a greedy manner where the KL divergence between distribution Q onto the convex hull of a set of distributions Q (composite null) is minimized. It is assumed that the KL divergence between Q and any distribution  $Q \in Q$  is finite  $^1$  (often call "nondegenerate" condition).

#### **Algorithm 1:** Li's Algorithm

```
\begin{array}{ll} \mathbf{1} & Q_{(1)} = \arg\min_{Q \in \mathcal{Q}} D(P\|Q) \\ \mathbf{2} & \textbf{for } m = 2, 3, \dots, K \textbf{ do} \\ \mathbf{3} & Q := \alpha Q_{(m-1)} + (1-\alpha)Q' \\ \mathbf{4} & \alpha, Q' \leftarrow \arg\min_{\alpha, Q'} D(P\|Q) \end{array}
```

Here, the distribution Q' and  $\alpha$  is chosen (coupled) such that the divergence is minimized. The minimizer need not be unique.

#### Regularity condition on alternative

Li's algorithm is apparently greedy with high fluctuation in initial steps. Additionally, this task is computationally expensive, and it is not clear of the convexity.

### Is the returned mixture in the convex hull?

The returned  $Q_{(m)}$  is in the convex hull. The first step returned a single element in Q with the smallest KL divergence. Iteratively, the linear combination is still in the convex hull, i.e.

$$\begin{split} Q_{(2)} &= \alpha_1 \cdot Q_{(1)} + (1 - \alpha_1) \cdot Q_{(1)}' \\ Q_{(2)} &= \alpha_2 \cdot Q_{(2)} + (1 - \alpha_2) \cdot Q_{(2)}' \\ &= \alpha_2 \cdot \alpha_1 \cdot Q_{(1)} + \alpha_2 \cdot (1 - \alpha_1) \cdot Q_{(1)}' + (1 - \alpha_2) \cdot Q_{(2)}'. \end{split}$$

It is clear that  $Q_{(2)}$  or  $Q_{(m)}$  would still be a convex combination of elements in Q.

4.2. Cisszar Algorithm. Originally proposed by

<sup>&</sup>lt;sup>1</sup>Which should I use? Q or ℚ

#### **5. Integrals.** What is it other than the area under curve?

Interesting examples where the Riemann's integration fails is that f(x) = 1 if x is rational and 0 otherwise over the interval [0,1]. Why this function's integral is problematic in Riemann's definition.

Lebesgue used measure theory to define integral.

Many times, when Riemann fails, Lebesgue works.

Set being countable and uncountable rely on if we can arrange them in a 'readable' sequence, could be infinite like rational numbers.

Proving that  $\mathbb{R}$  is uncountable is not so trivial: decimal expansion.

No matter how hard you try to write the numbers in [0,1), there is always some values left out.

Continuum hypothesis: How do you check which  $\infty$  is bigger? Or which set has the higher cardinality?

Definition 1.2.2 ()

EXAMPLE 5.1 (Set of Lebesgue measure zero, 1.2.2). A set S is zero in Lebesgue measure when it can be covered with a sequence of open intervals  $(I_1, I_2, \ldots)$ . And the sum  $\sum_{0}^{\infty} m(I_n)$  can be made arbitrarily small

In other words, for any  $\epsilon > 0$ , we can always find a sequence of  $I_n$  that covers S.

Theorem 5.2. Any countable infinite set of S has Lebesgue measure zero.

REMARK 5.3. The set of measure zero would not change a damn thing on f

Before writing this section, I sometimes came across expressions such as  $p(x) = A \, \mathrm{d} x$  or  $p(x) = A' m(\mathrm{d} x)$  in defining exponential family. It relates to the fundamental concept of measure and a density in measure theory. When I see  $\mathrm{d} x$  in evaluating integral, e.g.  $f(x) = \int_{\mathbb{R}} x \, \mathrm{d} x = x^2$ , it usually refers to the Lebesgue measure. In the case of  $m(\mathrm{d} x)$ , it generally emphasize that you are integrating with respect to a general measure m.

Wait, then what are we talking in defining distribution?

**6. Exponential Family.** The exponential family is a collection of parametric models with very elegant properties. Peter and Hao utilized it to arrive some nice results about the optimality of E-variables, mainly relying on the duality. We called a model belonging to the exponential family if the underlying distributions can be written below.

DEFINITION 6.1 (Exponential Family).

(6.1) 
$$p_{\beta}(u) := \exp\left(\beta^T t(u) - \psi(\beta)\right) p_0(u) m(\mathrm{d}u)$$

(6.2) 
$$= \frac{1}{Z(\beta)} \exp(\beta^T y) p_0(y) m(\mathrm{d}y).$$

- u: random variable,  $d \times 1$  vector in  $\mathcal{U} \subset \mathbb{R}^d$
- t(u): sufficient statistics,
- $\beta$ : canonical parameter,  $d \times 1$  vector in  $B \subset \mathbb{R}^d$
- $\psi(\beta)$ : cumulant function
- $p_0(y)$ : the *carrying density*, defined with respect to some *carrying measure* m(dy) on  $\mathcal{U}$ . We write the *carrying measure* explicitly m(du).
- $Z(\beta)$ : the partition function,  $Z(\beta) = \exp(\psi(\beta))$

Is the space for u and t(u) the same?

6.1. Basic Properties. Cumulant generating function  $\phi(\beta)$ : The first and second *cumulant* is mean and variance. Also  $\psi(\beta)$  is differentiable infinitely often (? true for all?).

DEFINITION 6.2 (Canonical Parameter Space). Canonical parameter space B is the set of parameter  $\beta$  where the following integration is finite

$$\mathtt{B} := \left\{ \beta : \int_{\mathcal{U}} \exp \left( \beta^T t(u) \right) p_0(u) m(\mathrm{d}u) < \infty \right\}.$$

 $m(\mathrm{d}u)$  will either be Lebesgue measures or counting measures for discrete cases. In the first case, ' $m(\mathrm{d}u)$ ' reduces to ' $\mathrm{d}u$ ', which can be handled using standard multivariable calculus. While in the latter, the above integral can be written as a summation.

REMARK 6.3. Every exponential family has a canonical parameterization with carrier density  $p_0$  such that the canonical parameter space contains the origin, i.e.  $0 \in B$ . (Section 18.4.3 in MDL)

PROOF. If  $0 \notin B$  or  $p_0$  is not a carrier density, we can pick any  $\beta_0 \in B$  and set

 $p'_0$ 

Now the  $\Box$ 

DEFINITION 6.4 (Regular Exponential Family). If the canonical parameter space of exponential families are nonempty open set, we call such families *regular*.

DEFINITION 6.5 (Full Exponential Family). The family is called *full* if the dimension of  $\beta$  equals the dimension of B.

DEFINITION 6.6 (Order). The order of an exponential family is the minimal dimension of t(u) such that we can express the family using Eq. (6.1).

DEFINITION 6.7 (Minimal Exponential Family). An exponential family is referred to as *minimal* if: a) there are no linear constraints among the components of the parameter vector  $\beta$ ; b) there are no linear constraints among the components of the sufficient statistic t(u) (in the latter case, with probability one under the measure m).

EXAMPLE 6.8 (Non-minimal distribution). The simplest would be a multinomial distribution with parameter (). The PMF can be written as

s

We can reparameterize the probability distribution even in the case of minimal distribution. Given a (?arbitrary) set  $\Theta$  and a mapping  $\Phi:\Theta\to B$ , we consider the densities with canonical parameters replaced by  $\Phi(\theta)$ 

$$p_{\theta}(u) := \exp(\Phi(\theta)^T t(u) - \phi(\Psi(\theta))).$$

Here  $\Phi$  is a one-to-one mapping whose image is all of B.

If  $\Phi$ 's image is a strict subset of B ( $\Phi(\Theta) \subset B$ ), it is then OK to reparameterize on that subset. If it can not be reducible, we refer this exponential family as curved.

We are also interested in cases in which the image of is a strict subset of N . If this subset is a linear subset, then it is possible to transform the representation into an exponential family on that subset. When the representation is not reducible in this way, we refer to the exponential family as a curved exponential family.

DEFINITION 6.9 (Curved Exponential Family). TBH

EXAMPLE 6.10 (Normal distribution with variance equalling to mean). Let  $u \in \mathbb{R} \sim \mathcal{N}(\mu, \mu^2)$ ,  $\mu \neq 0$ . The density of u is

$$p_{\mu}(u) = |\mu|^{-1} (2\pi)^{-1/2} \exp\left(-\frac{1}{2} (\frac{u-\mu}{\mu})^2\right)$$
$$= (2\pi)^{-1/2} \exp\left(-\frac{u^2}{2\mu^2} + \frac{u}{\mu} - \frac{1}{2} + \log(|\mu|)\right)$$
$$= \exp(\beta^T t(u) - \psi(\beta)).$$

The  $\beta^T = (-1/2\mu^2, 1/\mu)$ , the sufficient statistics  $t(u) = (u^2, u)$ . The dimension of the sufficient statistic is more than the dimension of  $\beta$  for curved exp. family.

I have found some other versions of the same statement regarding minimal exponential family:

- A minimal exponential family is where the t(u) are linearly independent
- A minimal exponential family is one where representation reaches the *order*

add example

All share the same statement regarding the linear dependency of sufficient statistics t(u), only Michael I. Jordan's Chapter 8 stated on  $\beta$ . I wonder if linear dependency in  $\beta$  implies linear dependency in t(u), i.e.

$$a^T \beta = C \Leftrightarrow b^T t(u) = C'$$

He also claimed that non-minimal families can always be reduced to minimal families via a suitable transformation and reparameterization.

If an exponential family is not minimal, it is called *overcomplete*. Both minimal and overcomplete representations are useful

- 6.2. Mean-value parameterization.
- 6.3. Simple Case.
- Why I think this is the carrier probability with some measure v  $p_{\mu^*} = p_{0,\mu^*}$   $q_{0,\mu^*} = q = q_{\mu^*}$ :

It is rather simple when you write down  $\beta = 0$ 

$$p_{\mathbf{0};\boldsymbol{\mu}^*}(u) := \frac{1}{Z_p(\mathbf{0};\boldsymbol{\mu}^*)} \exp(\mathbf{0}^T t(u)) \cdot p_{\boldsymbol{\mu}^*}(u)$$

$$= \frac{p_{\boldsymbol{\mu}^*}(u)}{Z_p(\mathbf{0};\boldsymbol{\mu}^*)}$$

$$= \frac{p_{\boldsymbol{\mu}^*}(u)}{\int \exp(\mathbf{0}^T t(u)) p_{\boldsymbol{\mu}^*}(u) \, \mathrm{d}\nu} = p_{\boldsymbol{\mu}^*}(u)$$

- A naive question follows: what is the distribution  $p_{\mu^*}(u)$ ? Is it in the mean-value parameterization or canonical form?
  - 6.4. General Case.
- Why Hao explicitly denote the mean of the carrier density in Grünwald, de Heide and Koolen (2024a)

We first discuss the Gaussian location family where the mean  $\mu^* \in M_p = \mathbb{R}^d$ . The null distribution will be family with  $\mu^*$  and a positive semidefinite covariance matrix  $\Sigma_p$  and the alternative  $\mathcal{Q} = \{Q\}$  would be Gaussian distribution with the same  $\boldsymbol{m}\boldsymbol{u}^*$  and a covariance matrix  $\Sigma_p \neq \Sigma_q$ .

The notation  $D_{\text{GAUSS}}$  in Eq. (3.2.1) in Hao (2025) is just the KL divergence. Here  $X = U = (X_1, \dots, X_d)$  is the d-dimensional random vector with distribution p or q

$$\begin{split} D_{\text{GAUSS}}(B) &= \text{D}_{\text{KL}}(P \parallel Q) \\ &:= \int_{\mathcal{U}} p(X) \log \frac{p(X)}{q(X)} = \mathbb{E}_p \left[ \log \frac{p(X)}{q(X)} \right] \\ &= \mathbb{E}_p \left[ \log \frac{(2\pi)^{-d/2} \det(\Sigma_p)^{-1/2} \exp\left(-\frac{1}{2}(X - \pmb{\mu}^*)^T \Sigma_p^{-1}(X - \pmb{\mu}^*)\right)}{(2\pi)^{-d/2} \det(\Sigma_q)^{-1/2} \exp\left(-\frac{1}{2}(X - \pmb{\mu}^*)^T \Sigma_q^{-1}(X - \pmb{\mu}^*)\right)} \right] \\ &= \mathbb{E}_p \left[ \log \frac{\det(\Sigma_q)^{-1/2} \exp\left(\Sigma_p^{-1}\right)}{\det(\Sigma_q)^{-1/2} \exp\left(\Sigma_q^{-1}\right)} \right] \quad \text{Wrong!} \end{split}$$

For more details, check https://statproofbook.github.io/P/mvn-kl.html for cases where the means are different.

Why Eq. (3.4.2) holds? The part I don't understand is that why putting a prior will equal to  $p_{\mu^*}$ 

$$S_{\text{COND}} := \frac{q_{W_1}(U^{(n)} \mid Z)}{p_{W_0}(U^{(n)} \mid Z)} \stackrel{?}{=} \frac{q_{\boldsymbol{\mu}^*}(U^{(n)} \mid Z)}{p_{\boldsymbol{\mu}^*}(U^{(n)} \mid Z)}$$

The key is just simple derivation with the fact that  $\hat{X}_{|n} := \sum X_i/n \sim \mathcal{N}(n\boldsymbol{\mu}^*, n\Sigma_q)$ . Setting  $Z = \hat{X}_{|n}/\sqrt{n} \sim \mathcal{N}(\sqrt{n}\boldsymbol{\mu}^*, \Sigma_q)$ , we have

$$\begin{split} q_{W_1}(X^{(n)} \mid Z) &:= \frac{q_{W_1}(X^{(n)}, Z)}{q_{W_1}(Z)} = \frac{q_{W_1}(X^{(n)})}{q_{W_1}(Z)} \stackrel{iid}{=} \frac{\prod_{i=1}^n q_{W_1}(X_i)}{q_{W_1}(Z)} \\ &= \frac{\prod (2\pi)^{-d/2} (\det \Sigma_q)^{-1} \exp\left(-1/2(X_i - \boldsymbol{\mu}^*)^\top \Sigma_q^{-1}(X_i - \boldsymbol{\mu}^*)\right)}{(2\pi)^{-d/2} (\det \Sigma_q)^{-1} \exp\left(-1/2(Z - \sqrt{n}\boldsymbol{\mu}^*)^\top \Sigma_q^{-1}(Z - \sqrt{n}\boldsymbol{\mu}^*)\right)} \\ &= \frac{\exp\left(\sum_{i=1}^n X_i^\top \Sigma_q^{-1} X_i - 2\sqrt{n}\boldsymbol{\mu}^{*\top} \Sigma_q^{-1} Z + n\boldsymbol{\mu}^{*\top} \Sigma_q^{-1} \boldsymbol{\mu}^*\right)}{\exp\left(Z^\top \Sigma_q^{-1} Z - 2\sqrt{n}\boldsymbol{\mu}^{*\top} \Sigma_q^{-1} Z + n\boldsymbol{\mu}^{*\top} \Sigma_q^{-1} \boldsymbol{\mu}^*\right)} \\ &= \exp\left(\sum_{i=1}^n (X_i - Z)^\top \Sigma_q^{-1}(X_i - Z)\right) \end{split}$$

The equality follows due to the joint distribution  $X^{(n)}$  and Z is simply  $X^{(n)}$ . For UI case, I don't see any sign of splitting? It looks like just ML prequential settings What makes the E-process-ness? th Why the  $S_{\text{SEO}} = S_{\text{SEO,RIP}}$  in Thm. 2?

6.5. Discussion with Peter at 21.05.2025. Papers discussed: PNAS and general case For the simple case (the Gaussian location family in 3.2.1 in Hao (2025)), my interpretation is as followed: If negative semidefinite  $\Sigma_q - \Sigma_p$ , then we reduced to the simple case where the RIPr E-variable is not only in  $\operatorname{conv}(\mathcal{P})$  (the convex hull of the null dist.) but rather a element of  $\mathcal{P}$  itself.

Otherwise, we can find a prior on  $\mathcal{P}$  with sharp variance  $\mathcal{N}(\mu^*, (\Sigma_q - \Sigma_p)/n)$ . The results in turn suggests that we cast a sharp prior on  $\mathcal{P}$  with the same mean. However, extra argument about the alternative Q, I think we move to the distributions from Q that shares the same sufficient statistics. This is still the part where I am having some doubts. I am still sure is that what does it mean to operate on a 'enlarged' or 'modified' alternative even if the alternative  $Q = \{Q\}$  being just Gaussian with same mean and a different variance.

Another point is in the Anti-Simple case, then  $S_{Q, \text{SEQ}, \text{RIP}}^{(n)} = 1$ . First of all, the superscript  $S_{Q, \text{SEQ}, \text{RIP}}^{(n)} = 1$ . First of all, the superscript says it is considering a sequential settings but  $S_{Q, \text{SEQ}, \text{RIP}}^{(n)} = 1$ . First of all, the superscript  $S_{Q, \text{SEQ}, \text{RIP}}^{(n)} = 1$ . First of all, the superscript says it is considering a sequential settings but  $S_{Q, \text{SEQ}, \text{RIP}}^{(n)} = 1$ . First of all, the superscript says it is considering a sequential settings but  $S_{Q, \text{SEQ}, \text{RIP}}^{(n)} = 1$ . First of all, the superscript says it is considering a sequential settings but  $S_{Q, \text{SEQ}, \text{RIP}}^{(n)} = 1$ . First of all, the superscript says it is considering a sequential settings but  $S_{Q, \text{SEQ}, \text{RIP}}^{(n)} = 1$ . First of all, the superscript says it is considering a sequential settings but  $S_{Q, \text{SEQ}, \text{RIP}}^{(n)} = 1$ . First of all, the superscript says it is considering a sequential settings but  $S_{Q, \text{SEQ}, \text{RIP}}^{(n)} = 1$ . First of all, the superscript says it is considering a sequential settings but  $S_{Q, \text{SEQ}, \text{RIP}}^{(n)} = 1$ . First of all, the superscript says it is considering a sequential setting setting says it is considering to  $S_{Q, \text{SEQ}, \text{RIP}}^{(n)} = 1$ . First of all, the superscript says it is considering a sequential setting says it is singular says it is considering a sequential setting says it is singular says it is considering a sequential setting says it is singular says it is s

**Data spilting in UI** In a sense, the spilting is done sequentially or done across n data points in contrast to the  $D_0$   $D_1$  in the original paper. Now we are talking about the UI in (3.2.14) I think! The classical on would be  $L = L_0(\hat{\theta}_1)/L_0(\hat{\theta}_0)$ 

$$S_{\breve{\boldsymbol{\mu}},\mathrm{UI}}^{(n)} = \frac{\prod_{i=1}^{n} q_{\breve{\boldsymbol{\mu}}|_{i-1}(U_{(i)})}}{p_{\breve{\boldsymbol{\mu}}|_{n}(U^{(n)})}}$$

We have a regular ML estimator likelihood in the denominator after looking at the whole sequence n. For the nominator, we have a product of the ML likelihood for each time i. In the end, we have effectively losing just  $U_{(1)}$  in calculating likelihood.

**Beyond NP** We discussed about the similarities among the loss function in the ERM scheme and in this paper. The ERM loss is always with respect to some data (empirical) while here we are concerning about a data-dependent loss?

The problem or rather common difficulty with p-value in NP paradigm is that why do we report p-value at all? What does it mean for a small p-value when  $p \ll \alpha$ ?

Peter proposed a alternative as in we should report E-value instead. I think also the notation of seeing  $\alpha$  as the error probability is somewhat challenged.

**Inspiration** In vaccination study, we see a extremely small p-value on null but we are not allowed to make *new* decision without setting up new hypothesises and new studies. In other words, if we looked at a promising data *post-hoc*, there is really not much we can do based on the original data. But with E-value, he argued that new decision can be based/conditioned on the data.

Is l in Eq. 1 Grünwald (2024) just a new  $\alpha$ ?

In this loss function  $L(\kappa, a)$ ,  $\kappa$  is the true state of nature that we don't know? We just assumed if the data is coming from null or alternative.

In all, I think the main idea is to change from Type-I error safe to Type-I risk safe. I think section 2's main take home message is that any maximally compatible decision rule must be a E-variable

This is cited from Micheal I Jordan's notes: https://people.eecs.berkeley.edu/~jordan/courses/260-spring10/other-readings/chapter8.pdf

6.6. One-dimensional special case.

$$\mathcal{G} := \{g_{\eta}(y), \eta \in A, \in \mathcal{Y}\}, \quad A \text{ and } \mathcal{Y} \in \mathbb{R}^{p}$$
$$g_{\eta}(y) := \exp(\eta y - \psi(\eta)) g_{0}(y) m(\mathrm{d}y)$$

REMARK 6.11 (Carrying density in p.3 efron). Any  $g_{\eta_0}(y) \in \mathcal{G}$  could be the carrier density and the members of  $\mathcal{G}$  are absolutely continuos with respect to each other, i.e. their null measure sets agree.

REMARK 6.12. Cumulant generating function for  $\psi(\eta)$  originates from the old techniques for finding expectations, variances and higher-order moments.

6.7. Moment Relationships. We can differentiate  $\exp(\psi(\eta)) = \int_{\mathcal{V}} \exp(\eta y) g_0(y) m(\mathrm{d}y)$ 

$$\dot{\psi}(\eta) \exp(\psi(\eta)) = \int_{\mathcal{Y}} y \exp(\eta y) g_0(y) m(\mathrm{d}y)$$
$$\mathbb{E}_{g_0}[y] = \dot{\psi}(\eta) = \int_{\mathcal{Y}} y \exp(\eta y - \psi(\eta)) g_0(y) m(\mathrm{d}y)$$

Differentiating  $\exp(\psi(\eta))$  twice gives:

$$(\ddot{\psi}(\eta) + \dot{\psi}(\eta)) \exp(\psi(\eta)) = \int_{\mathcal{Y}} y^2 \exp(\eta y) g_0(y) m(\mathrm{d}y)$$
$$\operatorname{Var}(y) = \ddot{\psi}(\eta) = \int_{\mathcal{Y}} (y^2 - y) \exp(\eta y - \psi(\eta)) g_0(y) m(\mathrm{d}y)$$

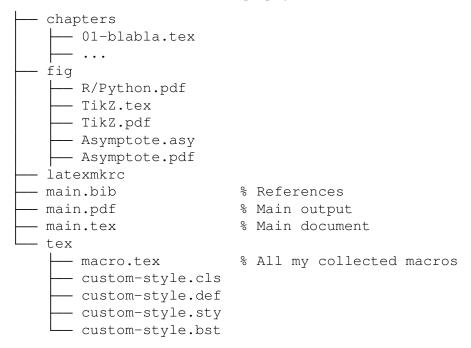
Efron's book (Chap. 2)

$$\mathcal{G} := \{g_{\eta}(y), \eta \in A, y \in \mathcal{Y}\}, \quad A \text{ and } \mathcal{Y} \in \mathbb{R}^{p}$$
$$g_{\eta}(y) := \exp\left(\eta^{T} y - \psi(\eta)\right) g_{0}(y) m(\mathrm{d}y)$$

- $g_0(y)$ : carrying density w.r.t. some carrying measure  $m(\mathrm{d}y)$  on  $\mathcal Y$  A is the canonical parameter space:
- $\psi(\eta)$  is the normalizing function or cumulant generating function (CGF)

**7. LATEX Project Management.** A consistent boilerplate for LATEX projects, I choose AOS for regular articles.https://vtex-soft.github.io/texsupport.ims-aos/

The references are managed externally by Zotero and BBT, exported to BibTex format, then included via natbib. Below is an example project hosted on Github or Overleaf:



- 7.1. Figure sizes and font sizes. Ratio, margin, font size, how to adjust accordingly...
- 7.2. Reference style.
- **8. Mathematical Notation.** It has always been a hassle to organise mathematical notation across different sources, in fact, I would go so far as to argue that it is the most annoying thing when one starts reading a book or an article.

However, there *must be* some notational conflicts beyond primary school simply due to the fact that the limited number of alphabets (26). For example, " $\mathbb{E}$ " might be energy in physics while it could refer to expectation or scores in probability.

Another difficulty is that the authors often assume some familiarity in the topics *also* I am expected to read in some logical or chronological order. In reality, I am constantly jumping back and forth between one literature to another.

$$\mathbb{A}, \mathbb{B}, \mathbb{C}, \mathbb{D}, \mathbb{E}, \mathbb{F}, \mathbb{G}, \mathbb{H}, \mathbb{J}, \mathbb{K}, \mathbb{L}, \mathbb{M}, \mathbb{N}, \mathbb{O}, \mathbb{P}, \mathbb{Q}, \mathbb{R}, \mathbb{S}, \mathbb{T}, \mathbb{U}, \mathbb{V}, \mathbb{W}, \mathbb{X}, \mathbb{Y}, \mathbb{Z}$$

$$\mathcal{A}, \mathcal{B}, \mathcal{C}, \mathcal{D}, \mathcal{E}, \mathcal{F}, \mathcal{G}, \mathcal{H}, \mathcal{I}, \mathcal{J}, \mathcal{K}, \mathcal{L}, \mathcal{M}, \mathcal{N}, \mathcal{O}, \mathcal{P}, \mathcal{Q}, \mathcal{R}, \mathcal{S}, \mathcal{T}, \mathcal{U}, \mathcal{V}, \mathcal{W}, \mathcal{X}, \mathcal{Y}, \mathcal{Z}$$

$$\mathfrak{A}, \mathfrak{B}, \mathfrak{C}, \mathfrak{D}, \mathfrak{E}, \mathfrak{F}, \mathfrak{G}, \mathfrak{H}, \mathfrak{I}, \mathfrak{I}, \mathfrak{I}, \mathfrak{K}, \mathfrak{L}, \mathfrak{M}, \mathfrak{N}, \mathfrak{O}, \mathfrak{P}, \mathfrak{Q}, \mathfrak{R}, \mathfrak{S}, \mathfrak{T}, \mathfrak{U}, \mathcal{V}, \mathcal{W}, \mathcal{X}, \mathcal{Y}, \mathcal{Z}$$

arg inf, arg sup, arg max, arg min, conv

This stackexchange answer <sup>2</sup> is probably the most comprehensive answer to which fonts are shown in LAT<sub>E</sub>X.

<sup>&</sup>lt;sup>2</sup>https://tex.stackexchange.com/a/58124

Symbol	Usage	Comments
$\mathbb{B}$	\*f	blackboard bold except \If due to conflict
${\cal B}$	\*C	calligraphic font
$\mathfrak{B}$	\*k	Fraktur font

8.1. Why the Fuck so many different notation. I am not even talking about the difference due to differences in fonts and italic or roman. It is just a very sad thing that we don't even have a unified way of saying probability is just sad.

Take probability for example and return to the most basic case of tossing a coin where the sample space is  $A = \{H, T\}$ . I have come across:

$$P()$$
  $P()$   $Pr()$   $P()$   $\mathbf{Prob}()$ .

From my experience, []

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