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# Fundamentals of Prequential Analysis

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*Prequential* = [Probabilistic]/Predictive/Sequential

— a general framework for assessing and comparing the predictive performance of a **FORECASTING SYSTEM**.

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- We assume reasonably extensive data, that either arrive in a time-ordered stream, or can be arranged into such a form:

$$\mathbf{X} = (X_1, X_2, \dots).$$

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- We try to identify these patterns, so as to use currently available data to form good forecasts of future values.

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**Basic idea:** Assess our future predictive performance by means of our past predictive performance.

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- Introduce the data-points  $(x_1, \dots, x_n)$  one by one.



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- Introduce the data-points  $(x_1, \dots, x_n)$  one by one.
- At time  $i$ , we have observed values  $\mathbf{x}^i$  of  $\mathbf{X}^i := (X_1, \dots, X_i)$ .

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We can assess forecast quality either in **absolute** terms, or by **comparison** of alternative sets of forecasts.

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$t$	1	2	3	...
$f$				
$x$				

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**Forecast type:** Pretty arbitrary: e.g.

- ☐ Point forecast
- ☐ Action
- ☐ Probability distribution



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**No peeping:** Forecast of  $X_{i+1}$  made before its value is observed — unbiased assessment

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Very general idea, e.g.:

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Very general idea, e.g.:

**No system:** e.g. day-by-day weather forecasts

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**Probability model:** Fully specified joint distribution  $P$  for  $\mathbf{X}$   
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**Statistical model:** Family  $\mathcal{P} = \{P_\theta\}$  of joint distributions for  $\mathbf{X}$

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□ forecast  $f_{i+1} = P^*(X_{i+1} \mid \mathbf{X}^i = \mathbf{x}^i)$ , where  $P^*$  is formed from  $\mathcal{P}$  by **somehow estimating/eliminating  $\theta$** , using the currently available data  $\mathbf{X}^i = \mathbf{x}^i$

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**Collection of models** e.g. forecast  $X_{i+1}$  using model that has performed best up to time  $i$

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—based on a statistical model  $\mathcal{P} = \{P_\theta\}$  for  $\mathbf{X}$ .

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**Plug-in forecasting system** Given the past data  $\mathbf{x}^i$ , construct some **estimate**  $\hat{\theta}_i$  of  $\theta$  (e.g., by maximum likelihood), and proceed as if this were the true value:

$$P_{i+1}^*(X_{i+1}) = P_{\hat{\theta}_i}(X_{i+1} \mid \mathbf{x}^i).$$

**NB:** This requires re-estimating  $\theta$  with each new observation!

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**Bayesian forecasting system (BFS)** Let  $\pi(\theta)$  be a prior density for  $\theta$ , and  $\pi_i(\theta)$  the posterior based on the past data  $\mathbf{x}^i$ . Use this to mix the various  $\theta$ -specific forecasts:

$$P_{i+1}^*(X_{i+1}) = \int P_\theta(X_{i+1} \mid \mathbf{x}^i) \pi_i(\theta) d\theta.$$

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**Other** e.g. fiducial predictive distribution, ...

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Gaussian process:  $X_i \sim \mathcal{N}(\mu, \sigma^2)$ ,  $\text{corr}(X_i, X_j) = \rho$



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MLEs:

$$\begin{aligned}\hat{\mu}_n &= \overline{X}_n && \xrightarrow{L} \mathcal{N}(0, \rho\sigma^2) \\ \hat{\sigma}_n^2 &= n^{-1} \sum_{i=1}^n (X_i - \overline{X}_n)^2 && \xrightarrow{p} (1 - \rho)\sigma^2 \\ \hat{\rho}_n &= 0\end{aligned}$$

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— not classically consistent.

But the estimated predictive distribution  $\hat{P}_{n+1} = \mathcal{N}(\hat{\mu}_n, \hat{\sigma}_n^2)$  **does** approximate the true predictive distribution  $P_{n+1}$ :  
normal with mean  $\bar{x}_n + (1 - \rho)(\mu - \bar{x}_n)/\{n\rho + (1 - \rho)\}$  and  
variance  $(1 - \rho)\sigma^2 + \sigma^2/\{n\rho + (1 - \rho)\}$ .

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Well-calibrated  
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The assessment of the quality of a forecasting system in the light of a sequence of observed outcomes should depend only on the forecasts it in fact delivered for that sequence

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— and not, for example, on how it might have behaved for other sequences.

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- ☐ Binary variables ( $X_i$ )
- ☐ Realized values ( $x_i$ )
- ☐ Emitted probability forecasts ( $p_i$ )

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- ☐ Binary variables ( $X_i$ )
- ☐ Realized values ( $x_i$ )
- ☐ Emitted probability forecasts ( $p_i$ )

Want (??) the ( $p_i$ ) and ( $x_i$ ) to be close “on average”:

$$\bar{x}_n - \bar{p}_n \rightarrow 0$$

where  $\bar{x}_n$  is the average of all the ( $x_i$ ) up to time  $n$ , *etc.*



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- ☐ Binary variables ( $X_i$ )
- ☐ Realized values ( $x_i$ )
- ☐ Emitted probability forecasts ( $p_i$ )

Want (??) the ( $p_i$ ) and ( $x_i$ ) to be close “on average”:

$$\bar{x}_n - \bar{p}_n \rightarrow 0$$

where  $\bar{x}_n$  is the average of all the ( $x_i$ ) up to time  $n$ , etc.

**Probability calibration:** Fix  $\pi \in [0, 1]$ , average over only those times  $i$  when  $p_i$  is “close to”  $\pi$ :

$$\bar{x}'_n - \pi \rightarrow 0$$

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Probability	0.4	0.6	0.3	0.2	0.6	0.3	0.4	0.5	0.6	0.2	0.6	0.4	0.3	0.5
Outcome	0	0	1	0	1	0	1	1	1	0	1	0	0	1

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Probability	0.4	0.6	0.3	0.2	0.6	0.3	0.4	0.5	0.6	0.2	0.6	0.4	0.3	0.5
Outcome	0	0	1	0	1	0	1	1	1	0	1	0	0	1

Probability	0.2	0.3	0.4	0.5	0.6
$p$					
Instances $n$	2	3	3	2	4
Successes	0	1	1	2	3
$r$					
Proportion	0	0.33	0.33	1	0.75
$\hat{p}$					

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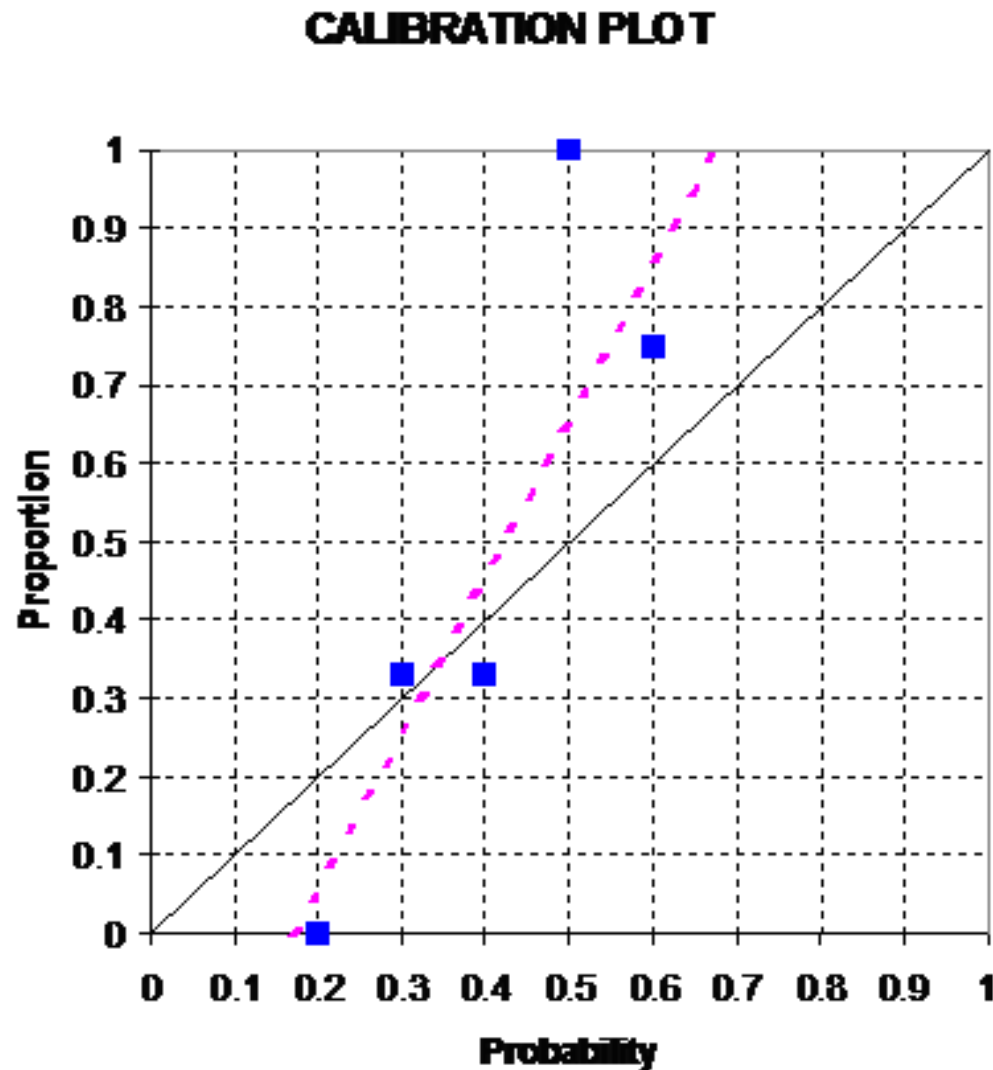
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Let  $\sigma$  be a **computable strategy** for selecting trials in the light of previous outcomes and forecasts

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Let  $\sigma$  be a **computable strategy** for selecting trials in the light of previous outcomes and forecasts

— e.g. third day following two successive rainy days, where forecast is below 0.5.

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Let  $\sigma$  be a **computable strategy** for selecting trials in the light of previous outcomes and forecasts

— e.g. third day following two successive rainy days, where forecast is below 0.5.

Then require asymptotic equality of averages,  $\bar{p}_\sigma$  and  $\bar{x}_\sigma$ , of the  $(p_i)$  and  $(x_i)$  over those trials selected by  $\sigma$ .

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**Why?**



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**Why?**

Can show following. Let  $P$  be a distribution for  $\mathbf{X}$ , and  $P_i := P(X_i = 1 \mid \mathbf{X}^{i-1})$ . Then

$$\bar{P}_\sigma - \bar{X}_\sigma \rightarrow 0$$

$P$ -almost surely, for **any** distribution  $P$ .

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Suppose  $p$  and  $q$  are computable forecast sequences, each computably calibrated for the same outcome sequence  $x$ .

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Suppose  $\mathbf{p}$  and  $\mathbf{q}$  are computable forecast sequences, each computably calibrated for the same outcome sequence  $\mathbf{x}$ .

Then  $p_i - q_i \rightarrow 0$ .

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Consider e.g.

$$Z_n := \frac{\sum (X_i - P_i)}{\sum P_i(1 - P_i)}$$

where  $P_i = P(X_i = 1 \mid X^{i-1})$ .

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Then

$$Z_n \xrightarrow{L} \mathcal{N}(0, 1)$$

for (almost) **any**  $P$ .

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So can refer value of  $Z_n$  to standard normal tables to test departure from calibration, **even without knowing generating distribution  $P$**

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— "Strong Prequential Principle"

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Suppose the  $X_i$  are continuous variables, and the forecast for  $X_i$  has the form of a continuous cumulative distribution function  $F_i(\cdot)$ .



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Suppose the  $X_i$  are continuous variables, and the forecast for  $X_i$  has the form of a continuous cumulative distribution function  $F_i(\cdot)$ .

If  $\mathbf{X} \sim P$ , and the forecasts are obtained from  $P$ :

$$F_i(x) := P(X_i \leq x \mid \mathbf{X}^{i-1} = \mathbf{x}^{i-1})$$

then, defining

$$U_i := F_i(X_i)$$

we have

$$U_i \sim U[0, 1]$$

independently, for any  $P$ .

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So we can apply various tests of uniformity and/or independence to the observed values

$$u_i := F_i(x_i)$$

to test the validity of the forecasts made

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So we can apply various tests of uniformity and/or independence to the observed values

$$u_i := F_i(x_i)$$

to test the validity of the forecasts made

— again, without needing to know the generating distribution  $P$ .

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Measure inadequacy of forecast  $f$  of outcome  $x$  by

loss function:  $L(x, f)$

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Measure inadequacy of forecast  $f$  of outcome  $x$  by

loss function:  $L(x, f)$

Then measure of overall inadequacy of forecast sequence  $\mathbf{f}^n$  for outcome sequence  $\mathbf{x}^n$  is **cumulative loss**:

$$L^n = \sum_{i=1}^n L(x_i, f_i)$$

We can use this to compare different forecasting systems.

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**Squared error:**  $f$  a point forecast of real-valued  $X$

$$L(x, f) = (x - f)^2.$$

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**Squared error:**  $f$  a point forecast of real-valued  $X$

$$L(x, f) = (x - f)^2.$$

**Logarithmic score:**  $f$  a probability density  $q(\cdot)$  for  $X$

$$L(x, q) = -\log q(x).$$



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At time  $i$ , having observed  $\mathbf{x}^i$ , probability forecast for  $X_{i+1}$  is its conditional distribution  $P_{i+1}(X_{i+1}) := P(X_{i+1} \mid \mathbf{X}^i = \mathbf{x}^i)$ .

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When we then observe  $X_{i+1} = x_{i+1}$ , the associated logarithmic score is

$$-\log p(x_{i+1} \mid \mathbf{x}^i).$$

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When we then observe  $X_{i+1} = x_{i+1}$ , the associated logarithmic score is

$$-\log p(x_{i+1} \mid \mathbf{x}^i).$$

So the cumulative score is

$$\begin{aligned} L_n(P) &= \sum_{i=0}^{n-1} -\log p(x_{i+1} \mid \mathbf{x}^i) \\ &= -\log \prod_{i=1}^n p(x_i \mid \mathbf{x}^{i-1}) \\ &= -\log p(\mathbf{x}^n) \end{aligned}$$

where  $p(\cdot)$  is the **joint density** of  $\mathbf{X}$  under  $P$ .

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$L_n(P)$  is just the (negative) **log-likelihood** of the joint distribution  $P$  for the observed data-sequence  $\mathbf{x}^n$ .

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$L_n(P)$  is just the (negative) **log-likelihood** of the joint distribution  $P$  for the observed data-sequence  $\mathbf{x}^n$ .

If  $P$  and  $Q$  are alternative joint distributions, considered as forecasting systems, then the excess score of  $Q$  over  $P$  is just the **log likelihood ratio** for comparing  $P$  to  $Q$  for the full data  $\mathbf{x}^n$ .

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This gives an interpretation to and use for likelihood that does not rely on the assuming the truth of any of the models considered.

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For a BFS:

$$\begin{aligned} P_{i+1}^*(X_{i+1}) &= \int P_{\theta}(X_{i+1} \mid \mathbf{x}^i) \pi_i(\theta) d\theta \\ &= P_B(X_{i+1} \mid \mathbf{x}^i) \end{aligned}$$

where  $P_B := \int P_{\theta} \pi(\theta) d\theta$  is the **Bayes mixture** joint distribution.

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where  $P_B := \int P_\theta \pi(\theta) d\theta$  is the **Bayes mixture** joint distribution.

This is equivalent to basing all forecasts on the single distribution  $P_B$ . The total logarithmic score is thus

$$\begin{aligned} L_n(\mathcal{P}) &= L_n(P_B) \\ &= -\log p_B(\mathbf{x}^n) \\ &= -\log \int p_\theta(\mathbf{x}^n) \pi(\theta) d\theta \end{aligned}$$



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For a plug-in system:  $L_n = -\log \prod_{i=0}^{n-1} p_{\hat{\theta}_i}(x_{i+1} \mid \mathbf{x}^i)$ .

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For a plug-in system:  $L_n = -\log \prod_{i=0}^{n-1} p_{\hat{\theta}_i}(x_{i+1} \mid \mathbf{x}^i)$ .

- The data  $(x_{i+1})$  used to evaluate performance, and the data  $(\mathbf{x}^i)$  used to estimate  $\theta$ , do not overlap
  - “unbiased” assessments (like cross-validation)

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  - “unbiased” assessments (like cross-validation)
- If  $x_i$  is used to forecast  $x_j$ , then  $x_j$  is *not* used to forecast  $x_i$ 
  - “uncorrelated” assessments (unlike cross-validation)

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- If  $x_i$  is used to forecast  $x_j$ , then  $x_j$  is *not* used to forecast  $x_i$ 
  - “uncorrelated” assessments (unlike cross-validation)

Both under- and over-fitting automatically and appropriately penalized.

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# Prequential efficiency

Let  $P$  be a SFS.  $P$  is prequentially **efficient** for  $\{P_\theta\}$  if, for **any** PFS  $Q$ :

$L_n(P) - L_n(Q)$  remains bounded above as  $n \rightarrow \infty$ ,  
with  $P_\theta$  probability 1, for almost all  $\theta$ .

[In particular, the losses of any two efficient SFS's differ by an amount that remains asymptotically bounded under almost all  $P_\theta$ .]

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[In particular, the losses of any two efficient SFS's differ by an amount that remains asymptotically bounded under almost all  $P_\theta$ .]

□ A BFS with  $\pi(\theta) > 0$  is prequentially efficient.

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Let  $P$  be a SFS.  $P$  is prequentially **efficient** for  $\{P_\theta\}$  if, for **any** PFS  $Q$ :

$L_n(P) - L_n(Q)$  remains bounded above as  $n \rightarrow \infty$ ,  
with  $P_\theta$  probability 1, for almost all  $\theta$ .

[In particular, the losses of any two efficient SFS's differ by an amount that remains asymptotically bounded under almost all  $P_\theta$ .]

- A BFS with  $\pi(\theta) > 0$  is prequentially efficient.
- A plug-in SFS based on a Fisher efficient estimator sequence is prequentially efficient.



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Model:

$$\mathbf{X} \sim P_{\theta} \quad (\theta \in \Theta)$$

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Model:

$$\mathbf{X} \sim P_{\theta} \quad (\theta \in \Theta)$$

Let  $P$  be prequentially efficient for  $\mathcal{P} = \{P_{\theta}\}$ , and define:

$$\begin{aligned}\mu_i &= \mathbb{E}_P(X_i \mid \mathbf{X}^{i-1}) \\ \sigma_i^2 &= \text{var}_P(X_i \mid \mathbf{X}^{i-1}) \\ Z_n &= \frac{\sum_{i=1}^n (X_i - \mu_i)}{\left(\sum_{i=1}^n \sigma_i^2\right)^{\frac{1}{2}}}\end{aligned}$$

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Then  $Z_n \xrightarrow{L} \mathcal{N}(0, 1)$  under **any**  $P_\theta \in \mathcal{P}$ .

So refer  $Z_n$  to standard normal tables to test the model  $\mathcal{P}$ .

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**Probability models** Collection  $\mathcal{C} = \{P_j : j = 1, 2, \dots\}$ .

- Both BFS and (suitable) plug-in SFS are prequentially **consistent**: with probability 1 under any  $P_j \in \mathcal{C}$ , their forecasts will come to agree with those made by  $P_j$ .

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**Parametric models** Collection  $\mathcal{C} = \{\mathcal{P}_j : j = 1, 2, \dots\}$ , where each  $\mathcal{P}_j$  is itself a parametric model:  $\mathcal{P}_j = \{P_{j,\theta_j}\}$ . Can have different dimensionalities.



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- Replace each  $\mathcal{P}_j$  by a prequentially efficient single distribution  $P_j$  and proceed as above.
- For each  $j$ , for almost all  $\theta_j$ , with probability 1 under  $P_{j,\theta_j}$  the resulting forecasts will come to agree with those made by  $P_{j,\theta_j}$ .

# Out-of-model performance

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Suppose we use a model  $\mathcal{P} = \{P_\theta\}$  for  $\mathbf{X}$ , but the data are generated from a distribution  $Q \notin \mathcal{P}$ . For an observed data-sequence  $\mathbf{x}$ , we have sequences of probability forecasts  $P_{\theta,i} := P_\theta(X_i \mid \mathbf{x}^{i-1})$ , based on each  $P_\theta \in \mathcal{P}$ : and “true” predictive distributions  $Q_i := Q(X_i \mid \mathbf{x}^{i-1})$ . The “best” value of  $\theta$ , for predicting  $\mathbf{x}^n$ , might be defined as:

$$\theta_n^Q := \arg \min_{\theta} \sum_{i=1}^n K(Q_i, P_{\theta,i}).$$

NB: This typically depends on the observed data

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With  $\hat{\theta}_n$  the maximum likelihood estimate based on  $\mathbf{x}^n$ , we can show that **for any  $Q$** , with  $Q$ -probability 1:

$$\hat{\theta}_n - \theta_n^Q \rightarrow 0.$$

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## Prequential analysis:

- ☐ is a natural approach to assessing and adjusting the empirical performance of a sequential forecasting system
- ☐ can allow for essentially arbitrary dependence across time
- ☐ has close connexions with Bayesian inference, stochastic complexity, penalized likelihood, *etc.*
- ☐ has many desirable theoretical properties, including automatic selection of the simplest model closest to that generating the data
- ☐ raises new computational challenges.

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Happy Birthday George!