

# STA 4103/5107 Computational Methods in Statistics II

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# **Special Topic 3**

# Laplace Approximation and Point Process Filter

**S3.1 Laplace Approximation** 



#### **Basic Idea**

- The idea behind the Laplace approximation is simple.
- We assume that the probability density  $P(x) = P^*(x)/Z_P$ , where the normalizing constant  $Z_P = \int P^*(x) dx$ .
- We also assume that the unnormalized density  $P^*(x)$  has a peak at a point  $x_0$ .
- We Taylor-expand the logarithm of  $P^*(x)$  around this peak:

where  $\log P^*(x) \approx \log P^*(x_0) - \frac{c}{2}(x - x_0)^2 + ...,$ 

$$c = -\frac{\partial^2}{\partial x^2} \log P^*(x) \Big|_{x=x_0}.$$



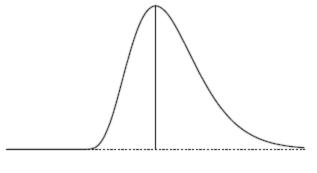
#### **Basic Idea**

• We then approximate  $P^*(x)$  by an unnormalized Gaussian,

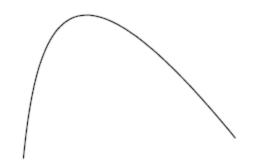
$$Q^*(x) \approx P^*(x_0) \exp[-\frac{c}{2}(x-x_0)^2],$$

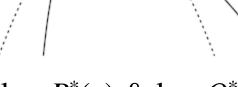
and we approximate the normalizing constant  $Z_P$  by the normalizing constant of this Gaussian,

$$Z_Q \approx \int P^*(x_0) \exp\left[-\frac{c}{2}(x - x_0)^2\right] dx = P^*(x_0) \sqrt{\frac{2\pi}{c}}$$









$$\log P^*(x) \& \log Q^*(x)$$



#### **General Case**

- We can generalize this integral to approximate  $Z_P$  for a density  $P^*(x)$  over a K-dimensional space x.
- If the matrix of second derivatives of  $\log P^*(x)$  at the maximum  $x_0$  is A, defined by:

$$A_{ij} = -\frac{\partial^2}{\partial x_i \partial x_j} \log P^*(x) \Big|_{x=x_0}.$$

Therefore,

$$\log P^*(x) \approx \log P^*(x_0) - \frac{1}{2} (x - x_0)^T A (x - x_0) + \dots,$$

• We let

$$Q^*(x) \approx P^*(x_0) \exp[-\frac{1}{2}(x-x_0)^T A(x-x_0)],$$



#### **General Case**

• Then, the normalizing constant is

$$Z_{Q} = \int P^{*}(x_{0}) \exp\left[-\frac{1}{2}(x - x_{0})^{T} A(x - x_{0})\right] dx$$
$$= P^{*}(x_{0}) \int \exp\left[-\frac{1}{2}(x - x_{0})^{T} A(x - x_{0})\right] dx$$

• Note that for k-dimensional random vector  $x \sim N(\mu, \Sigma)$ 

$$f(x) = \frac{1}{(2\pi)^{k/2} (\det \Sigma)^{1/2}} \exp\left[-\frac{1}{2} (x - \mu)^T \Sigma^{-1} (x - \mu)\right]$$

Therefore,

$$Z_Q = P^*(x_0) \sqrt{\frac{(2\pi)^k}{\det A}}.$$



# **Property**

- The Laplace approximation is a Gaussian based method.
- That is, the logarithm of a Gaussian function is a quadratic function, and this is the basic idea for the approximation.
- This method works well when the logarithm of the object function is concave.



#### **S3.2 Point Process Filter**



### **Filtering Estimation**

- Let  $x_1, x_2, \ldots, x_t, x_{t+1}, \ldots$ , form a process of interest, and instead of measuring  $x_t$ s directly, one observes variables  $y_1, y_2, \ldots, y_t$ ,  $y_{t+1}, \ldots$
- The goal is to use the observations, and joint probability models of x and y to estimate the unknown  $x_t$ s.
- Let  $f(x_t | y_1, y_2, ..., y_t)$  be the posterior density function of  $x_t$  given a set of observations  $(y_1, y_2, ..., y_t)$ .
- We are interested in the filtering estimation (posterior mean):

$$\hat{x}_t = E_f(x_t \mid y_1, y_2, ..., y_t).$$



# **Nonlinear Filtering Problem**

- Let the state vector  $x_t \in \mathbf{R}^d$  and the observation vector  $y_t \in \mathbf{R}^C$ .
- State equation:  $x_{t+1} = F(x_t) + w_t$ Observation equation:  $y_t = G(x_t) + q_t$
- Estimation:

update equation

$$f(x_t \mid y_1, y_2, ..., y_{t-1}) = \int_{x_{t-1}} f(x_t \mid x_{t-1}) f(x_{t-1} \mid y_1, y_2, ..., y_{t-1}) dx_{t-1}$$

prediction equation

$$f(x_t | y_1, y_2, ..., y_t) = \frac{f(y_t | x_t) f(x_t | y_1, y_2, ..., y_{t-1})}{f(y_t | y_1, y_2, ..., y_{t-1})}$$



#### **Point Process Observation**

- Assume the state vector  $x_k \in \mathbf{R}^d$
- Assume the observation is a **point process** in time interval [0, T].
- The time interval is discretized to time bins  $t_1, t_2, ..., t_M$ .
- For c = 1, ..., C, let the c-th observation component  $y_{k,c}$  = number of events in the k-th time bin.
- This number is either 0 or 1 if the bin size is sufficiently small.
- Therefore, the observation at the k-th time bin is

$$y_k = \{y_{k,c}\}_c \in \mathbf{R}^C$$



# **State-Space Model**

• Assume  $x_k$  follow a simple linear Gaussian transition. That is,

$$x_k = A x_{k-1} + w_k, \qquad w_k \in N(0, W)$$

where the A and W can be fitted in closed-form using **Maximum** likelihood Estimation (MLE).

• For  $y_k$ , we assume a generalized linear model (GLM) with an inhomogeneous Poisson process condition on  $x_k$ . That is,

$$y_{k,c} \sim Poisson(\lambda_{k,c})$$

where

$$\lambda_{k,c} = \exp(\mu_c + \alpha_c^T x_k)$$

•  $\{\mu_c, \alpha_c\}$  can also be identified using **MLE**.



# **System Identification**

- For each c = 1, ..., C, assume the observations are  $\{x_k, y_{k,c}\}$ .
- We maximize the likelihood

$$L = p(\{y_{k,c}\} | \{x_k\}) = \prod_{k=1}^{M} p(y_{k,c} | x_k) = \prod_{k=1}^{M} \frac{e^{-\lambda_{k,c}} (\lambda_{k,c})^{y_{k,c}}}{y_{k,c}!}$$

The log-likelihood is:

$$LL = \sum_{k=1}^{M} y_{k,c} \log(\lambda_{k,c}) - \lambda_{k,c} + const$$

$$= \sum_{k=1}^{M} y_{k,c} (\theta_c^T X_k) - \exp(\theta_c^T X_k) + const$$

where

$$\theta_c = (\mu_c, \alpha_c^T)^T, X_k = (1, x_k^T)^T$$



# **System Identification**

We use a Newton-Raphson method.

$$\frac{\partial LL}{\partial \theta_c} = \sum_{k=1}^{M} y_{k,c} X_k - \exp(\theta_c^T X_k) X_k$$

$$\frac{\partial^2 LL}{\partial \theta_c \partial \theta_c^T} = -\sum_{k=1}^{M} \exp(\theta_c^T X_k) X_k X_k^T$$

Recursive update:

$$(\theta_c)_{i+1} = (\theta_c)_i - \left(\frac{\partial^2 LL}{\partial \theta_c \partial \theta_c^T}\right)_i^{-1} \left(\frac{\partial LL}{\partial \theta_c}\right)_i$$



#### **Point Process Filter**

- To estimate the posterior  $f(x_k | y_1, y_2, ..., y_k)$ , we can use a sequential Monte Carlo (SMC) method.
- However, the method depends on number of sample points at each time step. (Note: these sample points are also called "particles" and the SMC is also called "particle filtering").
- A large number of particles often leads to inefficient computation.
- Here we introduce an efficient, deterministic estimation method, called point process filter.
- This method is based on Laplace approximation by approximate the posterior at each time using a Gaussian distribution.



#### **Estimation Process**

We use the recursive formula

$$f(x_k | y_1, y_2, ..., y_k) \propto f(y_k | x_k) f(x_k | y_1, y_2, ..., y_{k-1})$$

• Assume that conditioned on  $x_k$ , all components in  $y_k$  are independent. Therefore

$$f(y_k | x_k) = \prod_{c=1}^{C} f(y_{k,c} | x_k)$$

$$= \prod_{c=1}^{C} \exp(-\lambda_{k,c}) \frac{(\lambda_{k,c})^{y_{k,c}}}{y_{k,c}!}$$

We use the following notation to simplify the sub-index

$$a_{1:n} = a_1, a_2, ..., a_n.$$



#### **Time Update**

• We approximate the posterior using a Gaussian distribution at each time *k*. That is, let

$$x_{k|k} = E(x_k \mid y_{1:k})$$
  $W_{k|k} = Var(x_k \mid y_{1:k})$ 

Then

$$f(x_k \mid y_{1:k-1}) = \int f(x_k \mid x_{k-1}) f(x_{k-1} \mid y_{1:k-1}) dx_{k-1}$$

is also normally distributed.

The mean is computed as:

$$x_{k|k-1} = E(x_k \mid y_{1:k-1}) = E(Ax_{k-1} + w_k \mid y_{1:k-1})$$

$$= AE(x_{k-1} \mid y_{1:k-1})$$

$$= Ax_{k-1|k-1}$$



#### **Time Update**

The covariance is computed as:

$$W_{k|k-1} = Var(x_k \mid y_{1:k-1}) = Var(Ax_{k-1} + w_k \mid y_{1:k-1})$$

$$= Var(Ax_{k-1} \mid y_{1:k-1}) + Var(w_k)$$

$$= AW_{k-1|k-1}A^T + W$$

Therefore,

$$f(x_{k} | y_{1:k}) \propto f(y_{k} | x_{k}) f(x_{k} | y_{1:k-1})$$

$$= \left( \prod_{c=1}^{C} \exp(-\lambda_{k,c}) \frac{(\lambda_{k,c})^{y_{k,c}}}{y_{k,c}!} \right) \cdot \exp(-\frac{1}{2} (x_{k} - x_{k|k-1})^{T} W_{k|k-1}^{-1} (x_{k} - x_{k|k-1}))$$



### **Measurement Update**

Then, the logarithm of the posterior is

$$\log f(x_k \mid y_{1:k})$$

$$= \left(\sum_{c=1}^{C} y_{k,c} \log \lambda_{k,c} - \lambda_{k,c}\right) - \frac{1}{2} (x_k - x_{k|k-1})^T W_{k|k-1}^{-1} (x_k - x_{k|k-1}) + const$$

We approximate this posterior by a Gaussian distribution

$$\log f(x_k \mid y_{1:k}) = -\frac{1}{2} (x_k - x_{k|k})^T W_{k|k}^{-1} (x_k - x_{k|k}) + const$$

• Then,

$$\frac{1}{2}(x_k - x_{k|k})^T W_{k|k}^{-1}(x_k - x_{k|k})$$

$$= \frac{1}{2} (x_k - x_{k|k-1})^T W_{k|k-1}^{-1} (x_k - x_{k|k-1}) - \left( \sum_{c=1}^C y_{k,c} \log \lambda_{k,c} - \lambda_{k,c} \right) + const$$



### **Measurement Update**

• Differentiate w.r.t. to  $x_k$ , we have

$$\begin{aligned} W_{k|k}^{-1}(x_k - x_{k|k}) \\ &= W_{k|k-1}^{-1}(x_k - x_{k|k-1}) - \sum_{c=1}^{C} [y_{k,c} \frac{\partial \log \lambda_{k,c}}{\partial x_k} - \frac{\partial \lambda_{k,c}}{\partial x_k}] \\ &= W_{k|k-1}^{-1}(x_k - x_{k|k-1}) - \sum_{c=1}^{C} [y_{k,c} \alpha_c - \lambda_{k,c} \alpha_c] \end{aligned}$$

Differentiate again,

$$W_{k|k}^{-1} = W_{k|k-1}^{-1} + \sum_{c=1}^{C} \frac{\partial \lambda_{k,c}}{\partial x_k} \alpha_c = W_{k|k-1}^{-1} + \sum_{c=1}^{C} \alpha_c \lambda_{k,c} \alpha_c^T$$



### **Measurement Update**

• Let  $x_k = x_{k/k-1}$  after the second differentiation, we have

$$W_{k|k} = \left(W_{k|k-1}^{-1} + \sum_{c=1}^{C} \alpha_c \exp(\mu_c + \alpha_c^T x_{k|k-1}) \alpha_c^T\right)^{-1}$$

• Let  $x_k = x_{k/k-1}$  after the first differentiation, we have

$$W_{k|k}^{-1}(x_{k|k-1} - x_{k|k}) = -\sum_{c=1}^{C} [y_{k,c}\alpha_c - \exp(\mu + \alpha_c^T x_{k|k-1})\alpha_c]$$

Therefore,

$$x_{k|k} = x_{k|k-1} + W_{k|k} \sum_{c=1}^{C} [y_{k,c} - \exp(\mu_c + \alpha_c^T x_{k|k-1})] \alpha_c$$



# **Algorithm**

#### Point Process Filter Algorithm:

Update from time *k*-1 to *k*:

$$x_{k-1} \mid y_{1:k-1} \sim N(x_{k-1|k-1}, W_{k-1|k-1}) \rightarrow x_k \mid y_{1:k} \sim N(x_{k|k}, W_{k|k})$$

Time update: 
$$W_{k|k-1} = AW_{k-1|k-1}A^{T} + W$$
  
 $X_{k|k-1} = AX_{k-1|k-1}$ 

Measurement update:

$$W_{k|k} = \left(W_{k|k-1}^{-1} + \sum_{c=1}^{C} \alpha_c \exp(\mu_c + \alpha_c^T x_{k|k-1}) \alpha_c^T\right)^{-1}$$

$$x_{k|k} = x_{k|k-1} + W_{k|k} \sum_{c=1}^{C} [y_{k,c} - \exp(\mu_c + \alpha_c^T x_{k|k-1})] \alpha_c$$