0.1 Bayesian Analysis of Uniform distribution

Given
$$p(\theta) \sim Pa(b, K) = \begin{cases} \frac{Kb^k}{\theta^{K+1}} & \text{for } \theta \ge b \\ 0 & \text{otherwise} \end{cases}$$

Now with the above Pareto prior, the joint distribution of θ and $D = \{x_1, x_2, ..., x_N\}$ is:

$$p(D,\theta) = P(D|\theta)P(\theta) = \begin{cases} \frac{1}{\theta^N} \cdot \frac{Kb^K}{\theta^{K+1}} & \text{for } \theta \ge \max(\max(D), b) \\ 0 & \text{otherwise} \end{cases}$$

With the given definition of P(D) and taking m = max(D), the posterior $p(\theta|D)$ can now be defined as follows:

$$p(\theta|D) = \frac{P(\theta,D)}{P(D)} = \begin{cases} \frac{\frac{1}{\theta^N} \cdot \frac{Kb^K}{\theta^{K+1}}}{K} & \text{for } b \geq m \text{ and } b \leq \theta \\ \frac{\frac{1}{\theta^N} \cdot \frac{Kb^K}{\theta^{K+1}}}{(N+K)b^N} & \frac{1}{\theta^N} \cdot \frac{Kb^K}{\theta^{K+1}} \\ \frac{Kb^K}{(N+K)m^{N+K}} & \text{for } m > b \text{ and } m \leq \theta \end{cases}$$

$$0 \text{ otherwise}$$

(1)

(2)

This can be simplified and written as:

$$p(\theta|D) = \frac{P(\theta, D)}{P(D)} = \begin{cases} \frac{(N+K)b^{N+K}}{\theta^{N+K+1}} & \text{for } b \ge m \text{ and } b \le \theta \\ \frac{(N+K)m^{N+K}}{\theta^{N+K+1}} & \text{for } m > b \text{ and } m \le \theta \end{cases}$$

Note: For $b \ge m$, $p(\theta|D) = 0$ for $\theta < b$ and similarly for m > b $p(\theta|D) = 0$ for $\theta < m$. The posterior in fact has the same function functional form as a Pareto distribution: We know the Pareto distribution, Pa is defined as: $Pa(\alpha, n) = \frac{\alpha n^{\alpha}}{\theta^{\alpha+1}}$. If we compare this with the above distribution, we can see $\alpha \equiv (N + K)$ and $n \equiv \{m, b\}$

0.2 The Tramcar problem

0.2.1 Part A

Assumptions made:

• trams are numbered sequentially as integers starting from 0 to some upper bound θ . Thus the likelihood function p(x) can be defined as:

$$p(x) = f(x|\theta) = \begin{cases} \frac{1}{\theta} \text{ for } 0 \le x \le \theta \\ 0 \text{ otherwise} \end{cases}$$

• K=1 and b=1 and the D= $\{100\}$. This implies m= 100

Since m > b the posterior $p(\theta|D)$ is given as:

$$p(\theta|D) = \frac{(N+K)m^{N+K}}{\theta^{N+K+1}} = \frac{2.100^2}{\theta^3}$$

0.2.2 Part B

The posterior $p(\theta|D)$ as shown above is a Pareto distribution. Hence the mean and maximum posterior of $p(\theta|D)$ can derived using the properties of a Pareto distribution.

$$\mu(mean) = \frac{(N+K)m}{N+K-1} = \frac{2m}{1} = 200$$

$$MAP = mode = m = 100$$

0.2.3 Part C

The predictive density is given by $p(x|D) = \int_0^\infty p(x|\theta)p(\theta|D)d\theta$. As already stated in part A , $p(\theta|D) = 0$ for $\theta < m$ thus

$$p(x|D) = \int_{m}^{\infty} p(x|\theta)p(\theta|D)d\theta$$

$$p(x|D) = \int_{100}^{\infty} \frac{1}{\theta} \cdot \frac{2.100^{2}}{\theta^{3}} d\theta$$

$$= \left[\frac{-2.100^{2}}{3.\theta^{3}}\right]_{100}^{\infty}$$

$$= \frac{2}{3.100}$$

$$= \frac{1}{150}$$
(3)

0.2.4 Part D

From Part C, it is clear that the predictive distribution p(x|D) is a uniform distribution $U(x,\theta)$ where $\theta = 150$:

$$p(x|D) = \begin{cases} \frac{1}{150} \text{ for } 0 \le x \le 150\\ 0 \text{ otherwise} \end{cases}$$

Therefore for a new data point \mathbf{x} , the prediction is :

$$p(\mathbf{x}|D) = \frac{1}{150}I(\mathbf{x} \in [0, 150]) = \frac{1}{150}I(\mathbf{x} \le 150)$$

For observations whose value lie outside 150, the probability of observing them given the dataset D is 0. Thus $p(50|D) = \frac{1}{150}$ and p(500|D) = 0.

0.2.5 Part E

As $K \longrightarrow 0$

 $\lim_{K\to 0} p(\theta) = 0.$ Thus the limit of posterior $p(\theta|D)$ when K tends to 0 :

$$\lim_{K \to 0} p(\theta|D) = \frac{P(\theta, D)}{P(D)} = \begin{cases} \frac{(N)b^N}{\theta^{N+1}} \text{ for } b \ge m \text{ and } b \le \theta \\ \frac{(N)m^N}{\theta^{N+1}} \text{ for } m > b \text{ and } m \le \theta \end{cases}$$

Observations:

As K → 0, the posterior p(θ|D) at the limiting value of K is still a Pareto distribution.
 Sending K to 0 tends to change only the shape parameter of the original posterior distribution which is illustrated as follows:

We know $Pa(\kappa, n) = \frac{\kappa n^{\alpha}}{\theta^{\kappa+1}}$.

By comparing the above distribution with our original posterior $p(\theta|D)$, we concluded that $\kappa \equiv (N+K)$ but as $K \to 0$ $\kappa \equiv N$

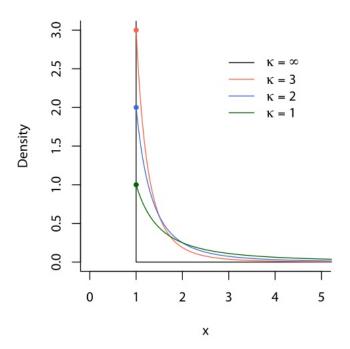
By observing the graph below (where x denote θ), we can see that decreasing the value of the shape parameter κ makes the distribution more heavily tailed. Thus, rare instances will have a greater probability mass assigned to them than before. Having the posterior $p(\theta|D)$ more heavily tailed improves the predictive density that we computed in Part C.

$$\lim_{k \to 0} p(x|D) = \int_{100}^{\infty} \frac{1}{\theta} \cdot \frac{1 \cdot 100^1}{\theta^2} d\theta = \frac{1}{200}$$

Therefore for a new data point \mathbf{x} , the prediction is :

$$p(\mathbf{x}|D) = \frac{1}{200}I(\mathbf{x} \in [0, 200]) = \frac{1}{200}I(\mathbf{x} \le 200)$$

Although the individual probability of observing each individual tram within the range [0-150] is less than before, by assigning probability masses on trams between 150 and 200, the predictor improves prediction by increasing the set of trams that it believes that we might see next. In other words the predictor moves towards the true uniform distribution.



0.3 Bayesian Classification

Given:

Initially expert1's knowledge regarding the probability of a food delivery being Soylent Red rather than Soylent Yellow is given by a beta distribution with parameters $(n_0, n_1) = (10, 10)$ while expert2's initial knowledge is best described by the beta distribution with parameters $(n_0, n_1) = (100, 20)$.

Hence expert 1's prior belief is captured by $p(f) = f^{(10)-1}(1-f)^{(10)-1}$ where f is probability of seeing Soylent Red. and expert 2's corresponding prior belief is given by $p(f) = f^{(100)-1}(1-f)^{(20)-1}$.

The prior distributions of both experts is shown in figure 1. Observation

• Since the true value of f is $\frac{3}{4}$, it can be seen that expert 2 initially has a better opinion. Reason: Without seeing any data, if both experts were asked to predict the probability Soylent Red, then on average both experts will chose a value of f that is equal to the mean of the prior distribution. The mean of expert 1's prior distribution is $\frac{1}{2}$ whereas for expert2 the mean of the beta distribution is $\frac{100}{120}$. The expected value of the prior distribution for expert2 is closer to the true value. Hence, judging from these values taken by the means it can be observed that expert 1 believes that on average he will see equal number of Soylent Reds and Soylent yellows whereas Expert2 believes on average he will see more Soylent reds than Soylent yellows.

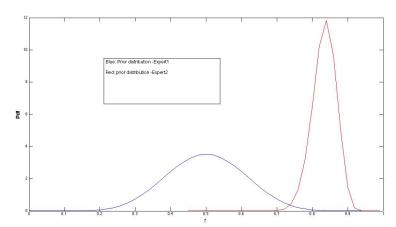


Figure 1: Prior beliefs of Expert1 and Expert2.

0.3.1Computing the posterior

With incoming data, the posterior distribution of both experts is given by:

$$p(f|\mathbf{X}) \propto p(\mathbf{X}|f)p(f)$$

$$= f^{k}(1-f)^{N-k}f^{n_{0}-1}(1-f)^{n_{1}-1}$$

$$= f^{(k+n_{0})-1}(1-f)^{(N+n_{1}-k)-1}$$
(4)

here k represents the number of food items seen that correspond to Soylent red and N corresponds to the total number of food items that has been seen.

For Expert 1, the corresponding posterior distribution is given a

$$p(f|X) = f^{(k+10)-1}(1-f)^{(N+10-k)-1}$$

And similarly, for expert 2 the corresponding posterior distribution $p(f|X) = f^{(k+100)-1}(1-f)^{(N+20-k)-1}$

$$p(f|X) = f^{(k+100)-1}(1-f)^{(N+20-k)-1}$$

Note: The mean and mode of the posterior distributions of both experts change as more observations made. The mean of the posterior distribution for expert 1 is $\mu_1 = \frac{K+10}{N+20}$ and

similarly the mean of the posterior distribution of expert 2 is $\mu_1 = \frac{K+100}{N+120}$. The terms K and N change with more data arrives. Hence, we can conclude that the belief of both experts change as they continue seeing new data. This is illustrated by the graph in figure 2.

Observation: Due to the high value for the n_0 parameter

Figure 2: Empirical mean over ${\bf k}$ for each posterior

