# Motivation for Fisher Information, Continued Again

## Seedlings (Poisson Model)

Ecologists divided a region of the forest floor into n quadrats and counted the number of seedlings that sprouted in each quadrat as part of a study on climate change.

- Observe  $X_1, \ldots, X_n$ ;  $X_i$  is the number of seedlings in quadrat number i.
- Data Model:  $X_i | \Lambda = \lambda$   $\stackrel{\text{i.i.d.}}{\sim} \text{Poisson}(\lambda)$  We have seen that the maximum likelihood estimate is  $\hat{\lambda}^{MLE} = \frac{1}{n} \sum_{i=1}^{n} X_i$

#### Connection between Observed Fisher Information and Taylor series approximation to log-likelihood

• Consider just the subset with 56 observations. The MLE is:

seedlings\new\_1993[subset\_inds]

## [36] 2 1 0 1 0 3 1 0 0 0 1 1 1 0 1 3 1 2 3 0 1

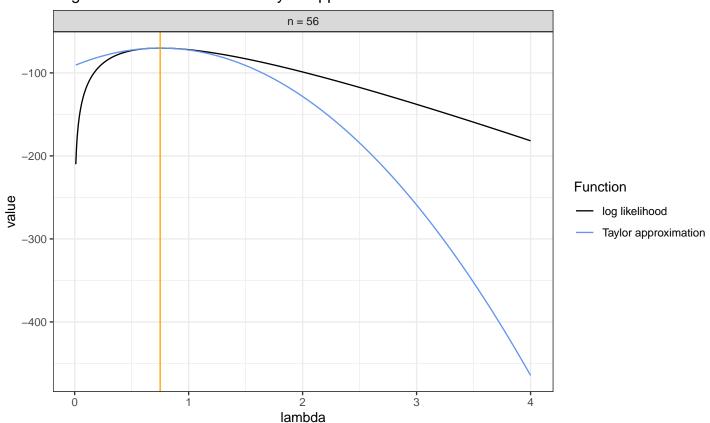
mean(seedlings\$new\_1993[subset\_inds])

## [1] 0.75

- There are n = 56 observations
- The observed Fisher information is  $J(\theta^*) = \frac{n}{\bar{x}} = \frac{56}{0.75} = 74.667$ .

   This is the negative second derivative of the log likelihood function.
- The Taylor series approximation about the maximum likelihood estimate  $\hat{\lambda}^{MLE}$  is:
  - $-\ell(\lambda|x_1,\ldots,x_n) \approx \ell(0.75|x_1,\ldots,x_{56}) \frac{1}{2}74.667(\lambda 0.75)^2$
- Here is the log-likelihood function with an orange line at the MLE and the Taylor approximation in blue:

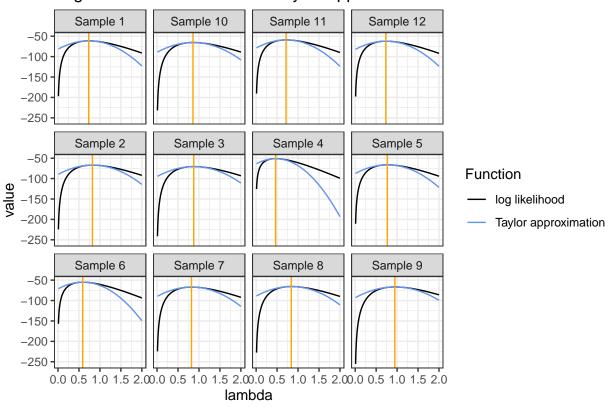
#### Log-likelihood Function and Taylor Approximation



#### What if we took other samples?

- Suppose for the sake of the example that the true parameter value is  $\lambda = 0.75$ .
- Each random sample of size n = 56 has:
  - Different observed values
  - A different log-likelihood function
  - A different second derivative of the log-likelihood function at the maximum
  - A different observed Fisher information

### Log-likelihood Function and Taylor Approximation



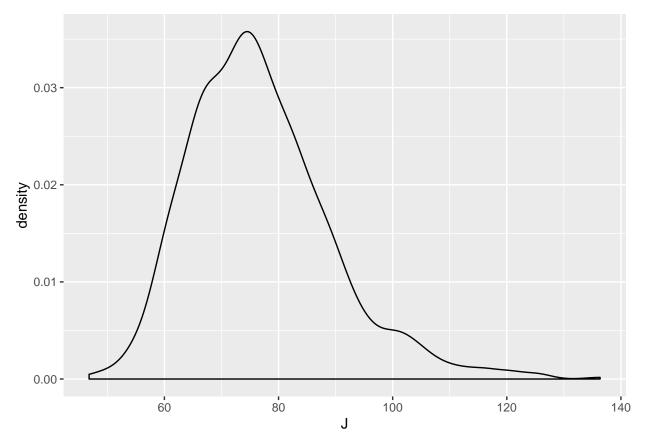
Suppose now I simulate 1000 different samples and calculate the observed Fisher information from each:

```
Fisher_informations <- data.frame(
    J = rep(NA, 1000)
)

for(i in 1:1000) {
    x <- rpois(n = 56, lambda = 0.75)
      Fisher_informations$J[i] <- 56 / mean(x)
}
head(Fisher_informations$J)

## [1] 65.33333 71.27273 78.40000 104.53333 69.68889 82.52632
length(Fisher_informations$J)

## [1] 1000
ggplot(data = Fisher_informations, mapping = aes(x = J)) +
    geom_density()</pre>
```



The Fisher information is the expected value (average) of the observed Fisher information across different samples: mean(Fisher\_informations\$J)

## [1] 76.76499