

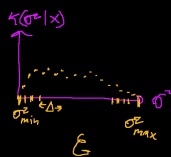
Grid Sampling Algorithm

1) Create a grid by first picking $\sigma_{\min}^2, \sigma_{\max}^2, \Delta$.

$$\mathcal{G} := \left\{ \sigma_{\min}^2, \sigma_{\min}^2 + \Delta, \sigma_{\min}^2 + 2\Delta, \dots, \sigma_{\max}^2 - \Delta, \sigma_{\max}^2 \right\}$$

For example, in our case let $\sigma_{\min}^2 = 0, \sigma_{\max}^2 = 1000, \Delta = 0.1$

$$\mathcal{G} = \{0, 0.1, 0.2, \dots, 999.9, 1000\}$$

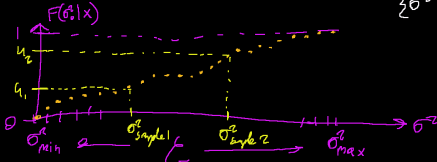


2) Approximate the value of c , the normalization constant via Riemann sum (if continuous).

$$c = \frac{1}{\int_{\text{supp}(\sigma^2)} k(\sigma^2|x) d\sigma^2} \approx \frac{1}{\Delta \sum_{\sigma^2 \in \mathcal{G}} k(\sigma^2|x)} = \frac{1}{\int_{-\infty}^{\infty} P(\sigma^2|x) d\sigma^2}$$

3) Compute the "sampling CDF" for all grid points:

$$F(\sigma_o^2|x) := P(\sigma^2 \leq \sigma_o^2 | x) \approx \sum_{\{\sigma^2: \sigma^2 \in \mathcal{G}, \sigma^2 \leq \sigma_o^2\}} c k(\sigma^2|x)$$



4) Draw u from $U(0,1)$ and locate $\sigma_{\text{sample}}^2 = \min_{\sigma^2 \in \mathcal{G}} \left\{ F(\sigma^2|x) \geq u \right\}$

If you want to draw many samples, repeat step 4. You only need to do steps 1-3 once.

Returning to our problem, how do we sample from the posterior $P(\theta, \sigma^2 | X)$?

- (1) Sample σ^2_{sample} from $P(\sigma^2 | X)$ using the grid sampler.
- (2) Sample θ_{sample} from $P(\theta | \sigma^2, X)$ via `rnorm`.

Return $[\theta_{\text{sample}}, \sigma^2_{\text{sample}}]$

At this point, we're not limited to distributions we know about nor limited to using conjugate priors. We're liberated!

We agree we can sample now from any arbitrary posterior. But how do we get point estimates?



I now have as many samples as I wish from my posterior (you want more? Let the computer churn more). You have S samples:

$$\begin{bmatrix} \theta_1 \\ \sigma_1^2 \end{bmatrix}, \begin{bmatrix} \theta_2 \\ \sigma_2^2 \end{bmatrix}, \dots, \begin{bmatrix} \theta_S \\ \sigma_S^2 \end{bmatrix}$$

How do we use these samples to get point estimates?

$$\hat{\theta}_{\text{MMSE}} \approx \frac{1}{S} \sum_{k=1}^S \theta_k$$

$\hat{\theta}_{\text{MAE}} \approx \text{SampleMedian}[\theta_1, \dots, \theta_S]$ i.e. you order all the samples and return the middle value

$\hat{\theta}_{\text{MAP}}$ complicated... you need you use smoothing... so forget it

$$\hat{\sigma}_{\text{MMSE}}^2 = \frac{1}{S} \sum_{k=1}^S \sigma_k^2$$

$$\hat{\sigma}_{\text{MAE}}^2 = \text{SampleMedian}[\sigma_1^2, \dots, \sigma_S^2]$$

How do we get CR's?

$$CR_{\theta, 1-\alpha_o} := [Q[\theta|x, \frac{\alpha_o}{2}], Q[\theta|x, 1-\frac{\alpha_o}{2}]]$$

$$\approx [\text{SampleQuantile}[\theta_1, \dots, \theta_S, \frac{\alpha_o}{2}], \text{SampleQuantile}[\theta_1, \dots, \theta_S, 1-\frac{\alpha_o}{2}]]$$

How do we get pvals for hypothesis tests?

$$H_0: \theta \in \Theta_0, \text{ pval} := P(\theta \in \Theta_0 | x) = \int_{\Theta_0} P(\theta|x) d\theta \approx \frac{1}{S} \sum_{k=1}^S \mathbb{1}_{\theta_k \in \Theta_0}$$

$$\text{e.g. } H_0: \theta \leq 5.89 \quad \text{pval} = \frac{\#\theta_k \leq 5.89}{S} = \text{proportion} \leq 5.89$$

How do we get the marginal distribution $P(\theta | X)$?

$$P(\theta|x) = \int_{\text{supp}(\sigma^2)} P(\theta, \sigma^2|x) d\sigma^2 \approx U(\underbrace{\{\theta_1, \dots, \theta_S\}}_{\text{samples}})$$

uniform discrete distribution

How do we get the posterior predictive distribution $P(X^* | X)$?

$$P(x_*|x) = \iint_{\mathbb{R} \times \mathbb{R}} \underbrace{P(x_*|\theta, \sigma^2)}_{N(\theta, \sigma^2)} \underbrace{P(\theta, \sigma^2|x)}_{??} d\theta d\sigma^2 = \iint P(x_*, \theta, \sigma^2|x) d\theta d\sigma^2$$

You use the sample $[\theta_1, \sigma^2_1]$ and then draw x^*_1 from `rnorm(θ_1, σ^2_1)`. Then use the sample $[\theta_2, \sigma^2_2]$ and then draw x^*_2 from `rnorm(θ_2, σ^2_2)` ..., Then use the sample $[\theta_S, \sigma^2_S]$ and then draw x^*_S from `rnorm(θ_S, σ^2_S)`:

$$U(\{x_{*1}, x_{*2}, \dots, x_{*S}\}) \approx P(x_*|x)$$

$$E[x_*|x] \approx \frac{1}{S} \sum_{k=1}^S x_{*k}$$

Disadvantages of the Grid Sampling Algorithm

- 1) In many dimensions, how do you pick min, max? For parameters in multiparameter models, it's not so simple!
- 2) Computers have numerical underflow and overflow
- 3) "Curse of dimensionality". Let's say you want to sample 10,000 grid points per dimension. But you have 10 different thetas. $10,000^{10} = 10^{50}$ which is IMPOSSIBLE for the computer. As another example: let's say I want 1 billion points in 10 dimensions. That is only: $\sqrt[10]{10^9} \approx 8$ points per dimension.

That's not good resolution at all!!!

We need another solution to this problem!!!