## Model Selection — Bayesian Information Criterion

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## Viewpoint from Homo-Bayesianis

The backdrop: We have a bunch of alternative models:  $\mathcal{M}_i$ , and each model gives a parameter space  $\Theta$  and a setting for generation of data  $p(\mathcal{D}|\boldsymbol{\theta}_i, \mathcal{M}_i)$ . For comparing fidelity of different models, under Bayesian principle, we should use  $p(\mathcal{D}|\mathcal{M}_i) \propto p(\mathcal{M}_i|\mathcal{D})$ , assuming the prior for different  $\mathcal{M}_i$  are equal. The previous  $p(\mathcal{M}_i|\mathcal{D})$  is called *model evidence*. By Bayes' formula, the evidence is obtained by:

$$p(\mathcal{D}|\mathcal{M}_i) = \int p(\mathcal{D}|\boldsymbol{\theta}_i, \mathcal{M}_i) \pi(\boldsymbol{\theta}_i|\mathcal{M}_i)) d\boldsymbol{\theta}$$
$$=: \int f_{\mathcal{M}_i; \mathcal{D}}(\boldsymbol{\theta}_i) d\boldsymbol{\theta}$$

So if the integration is hard to compute, it's reasonable to assume that  $f_{\mathcal{M}_i;\mathcal{D}}(\boldsymbol{\theta}_i)$  as a function of  $\boldsymbol{\theta}_i$  is close to a p.d.f. of a normal distribution [a homo-frequentitus will explain it by asymptotic normality], so by Laplace approximation near the MAP point  $\hat{\boldsymbol{\theta}}_i$ , the integration is approximately decided by the Hessian matrix of  $log\ f_{\mathcal{M}_i;\mathcal{D}}$  at the MAP, as follows:

$$log \ p(\mathcal{D}|\mathcal{M}_i) = log \ p(\mathcal{D}|\hat{\boldsymbol{\theta}}_i, \mathcal{M}_i) + log \ \pi(\hat{\boldsymbol{\theta}}_i|\mathcal{M}_i) + \frac{k}{2}log(2\pi) - \frac{1}{2}log|\mathbf{A}|$$

where k represents the dimension of the i-th parameter space and A is the negative Hessian matrix at  $\hat{\theta}_i$ :

$$\mathbf{A} = -\nabla^2 \left| \log p(\mathcal{D}|\boldsymbol{\theta}_i, \mathcal{M}_i) \pi(\boldsymbol{\theta}_i|\mathcal{M}_i) \right|$$

$$= -\nabla^2 \left| \log p(\mathcal{D}|\boldsymbol{\theta}_i, \mathcal{M}_i) + \log \pi(\boldsymbol{\theta}_i|\mathcal{M}_i) \right|$$

The last 3 terms of RHS are comprised as the penalization term against complexity, called "Occam factor", while the first term of RHS describes how well can a prediction under the model fit the given data. As N increases far larger than k: the first 2 terms of Occam factor  $\log \pi(\hat{\theta}_i|\mathcal{M}_i) + \frac{k}{2}\log(2\pi)$  can be relatively ignored and, for the last term:

$$\begin{aligned} |det(\mathbf{A})|^{\frac{1}{2}} \sim & \left| -\nabla^{2}|_{\hat{\boldsymbol{\theta}_{i}}}log \ p(\mathcal{D}|\boldsymbol{\theta}_{i}, \mathcal{M}_{i}) \right|^{\frac{1}{2}} \\ \sim & \left| -\nabla^{2}|_{\hat{\boldsymbol{\theta}_{i}}}N \cdot log \ p(x_{typical}|\boldsymbol{\theta}_{i}, \mathcal{M}_{i}) \right|^{\frac{1}{2}} \\ \sim & O(N^{\frac{k}{2}}) \end{aligned}$$

So we can approximately estimate  $log p(\mathcal{D}|\mathcal{M}_i)$  by

$$log \ p(\mathcal{D}|\hat{\boldsymbol{\theta}_i}, \mathcal{M}_i) - rac{k}{2}logN$$

which we should maximize among all models.

The BIC (which we should minimize, like AIC) is formally defined as

$$BIC = K \log N - 2 \log p(\mathcal{D}|\hat{\boldsymbol{\theta}}_i, \mathcal{M}_i)$$