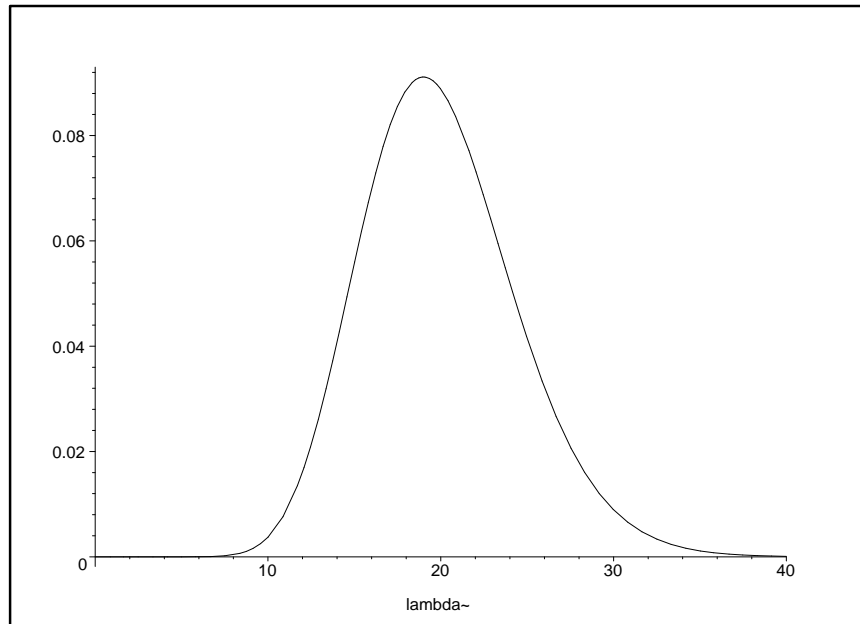


## Diffuse Priors (continued)

```
> plot( p( lambda, 20, 1 ), lambda = 0 .. 40, color = black );
```



$\Gamma(20, 1)$  does indeed look **not far from Gaussian**, and at first it may appear that it is indeed **relatively flat** in the region where the likelihood is appreciable ( $\lambda \in (1.0, 3.3)$ ), but we'll see below that it's actually **rather more informative** than we intend.

Recalling that the **mean** and **SD** of a  $\Gamma(\alpha, \beta)$  random quantity are  $\frac{\alpha}{\beta}$  and  $\sqrt{\frac{\alpha}{\beta^2}}$ , respectively, and that when used as a prior with the Poisson likelihood the  $\Gamma(\alpha, \beta)$  distribution acts like a dataset with **prior sample size**  $\beta$ , you can construct the following table:

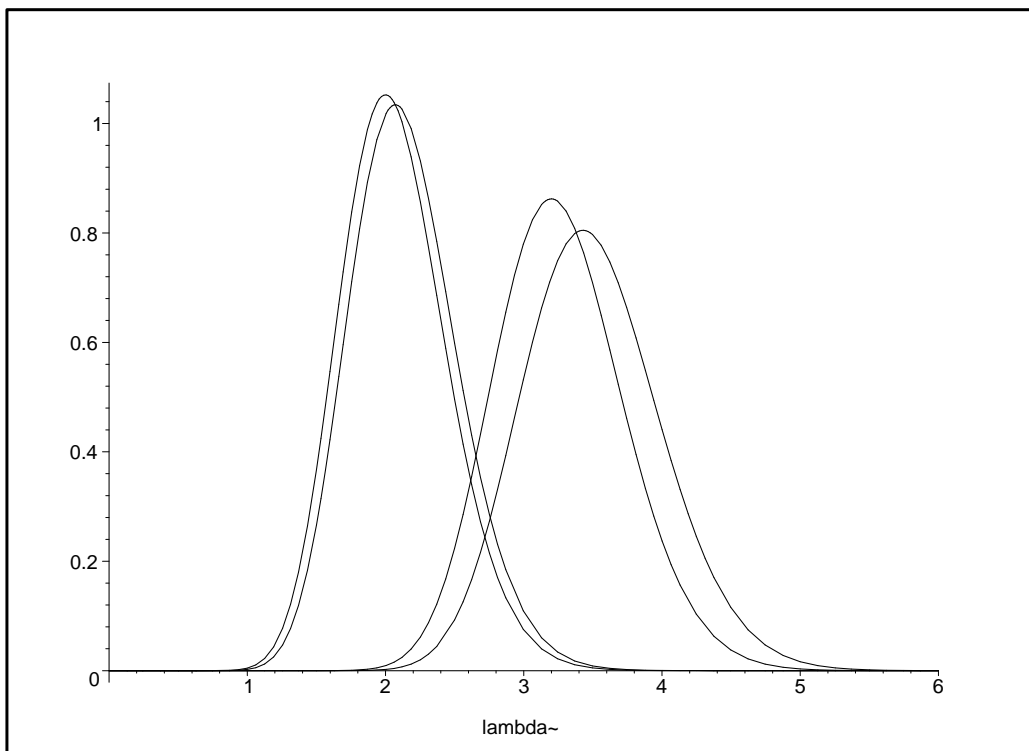
Prior				Posterior			
$\alpha$	$\beta =$ Sample Size	Mean	SD	$\alpha^*$	$\beta^*$	Mean	SD
0.001	0.001	1	31.6	29.001	14.001	2.071	0.385
1	0.001	1000	1000	30	14.001	2.143	0.391
20	1	20	4.47	49	15	3.267	0.467
20	0.001	20000	4472	49	14.001	3.500	0.500
$U(0, C)$ for $C > 4$		$\frac{C}{2}$	$\frac{C}{\sqrt{12}}$	30	14	2.143	0.391

## Diffuse Priors (continued)

The  $\Gamma(1, \epsilon)$  prior leads to an analysis that's **essentially equivalent** to the **integrated likelihood (fiducial)** approach back on p. 72, and the  $U(0, C)$  prior for  $C > 4$  (say) produces similar results:  $U(0, C)$  yields the  $\Gamma(s + 1, n)$  posterior **truncated** to the right of  $C$  (and this truncation has **no effect** if you choose  $C$  big enough).

You might say that the  $U(0, C)$  distribution has a **prior sample size of 0** in this analysis, and its prior mean  $\frac{C}{2}$  and SD  $\frac{C}{\sqrt{12}}$  (both of which can be made arbitrarily large by letting  $C$  grow without bound) are **irrelevant** (an example of how intuition can change when you depart from the class of **conjugate** priors).

```
> plot( { p( lambda, 29.001, 14.001 ), p( lambda, 30, 14.001 ),
        p( lambda, 49, 15 ), p( lambda, 49, 14.001 ) }, lambda = 0 .. 6,
        color = black );
```



The **moral** is that with only  $n = 14$  observations, some care is needed (e.g., through **pre-posterior** analysis) to achieve a prior that **doesn't affect the posterior very much**, if that's your goal.

## 2.8 Continuous Outcomes

For **continuous outcomes** there's an analogue of de Finetti's Theorem that's **equally central** to Bayesian model-building (e.g., Bernardo and Smith, 1994):

### de Finetti's Theorem for Continuous Outcomes.

If  $Y_1, Y_2, \dots$  is an infinitely exchangeable sequence of **real-valued** random quantities with probability measure  $p$ , there exists a probability measure  $Q$  over  $\mathcal{D}$ , the space of all distribution functions on  $R$ , such that the joint distribution function of  $Y_1, \dots, Y_n$  has the form

$$p(y_1, \dots, y_n) = \int_{\mathcal{D}} \prod_{i=1}^n F(y_i) dQ(F), \quad (77)$$

where  $Q(F) \stackrel{P}{=} \lim_{n \rightarrow \infty} p(F_n)$  and  $F_n$  is the **empirical distribution function** based on  $Y_1, \dots, Y_n$ .

In other words, exchangeability of real-valued observables is **equivalent** to the hierarchical model

$$\begin{array}{ccc} F & \sim & p(F) \quad (\text{prior}) \\ (Y_1, \dots, Y_n | F) & \stackrel{\text{IID}}{\sim} & F \quad (\text{likelihood}) \end{array} \quad (78)$$

for some **prior distribution**  $p$  on the **set**  $\mathcal{D}$  of all possible **distribution functions**.

This prior makes the continuous form of de Finetti's Theorem **considerably harder to apply**: to take the elicitation task seriously is to try to specify a probability distribution on a **function space** ( $F$  is in effect an **infinite-dimensional** parameter).

(**NB** This task is not unique to Bayesians—it's equivalent to asking "**Where does the likelihood come from?**" in frequentist analyses of observational data.)

# Continuous Outcomes (continued)

What people often do in practice is to appeal to considerations that narrow down the field, such as an *a priori* judgment that the  $Y_i$  ought to be **symmetrically** distributed about a measure of center  $\mu$ , and then try to use a fairly **rich parametric family** satisfying (e.g.) the symmetry restriction as a substitute for all of  $\mathcal{D}$ .

Strictly speaking you're not supposed to look at the  $Y_i$  while specifying your prior on  $\mathcal{D}$ —this can lead to a failure to fully assess and propagate **model uncertainty**—but not doing so can permit the data to surprise you in ways that would make you want to go back and revise your prior (an example of **Cromwell's Rule** in action).

As mentioned earlier, I'll suggest two potential ways out of this dilemma, based on **out-of-sample predictive validation** (the model-checking in the LOS data above was an example of this) and **Bayesian nonparametrics**.

**Case Study:** *Measurement of physical constants.* What used to be called the National Bureau of Standards (NBS) in Washington, DC, conducts extremely high precision measurement of physical constants, such as the actual weight of so-called **check-weights** that are supposed to serve as reference standards (like the official kg).

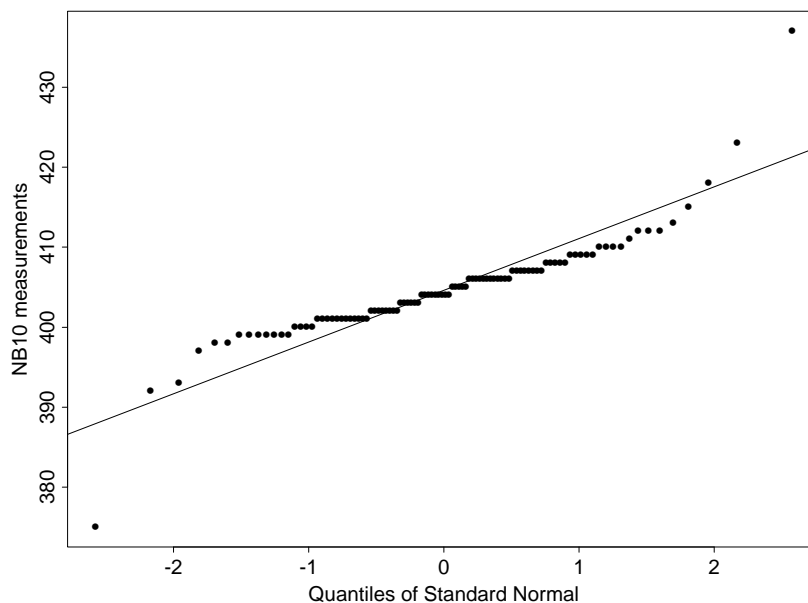
In 1962–63, for example,  $n = 100$  weighings (listed below) of a block of metal called **NB10**, which was supposed to weigh exactly 10g, were made under conditions **as close to IID as possible** (Freedman et al., 1998).

Value	375	392	393	397	398	399	400	401
Frequency	1	1	1	1	2	7	4	12
Value	402	403	404	405	406	407	408	409
Frequency	8	6	9	5	12	8	5	5
Value	410	411	412	413	415	418	423	437
Frequency	4	1	3	1	1	1	1	1

# NB10 Modeling

**Q:** (a) How much does NB10 **really weigh**? (b) How certain are you given the data that the true weight of NB10 is **less than** (say) 405.25? And (c) How accurately can you **predict** the 101st measurement?

The graph below is a **normal qqplot** of the 100 measurements  $y = (y_1, \dots, y_n)$ , which have a mean of  $\bar{y} = 404.6$  (the units are **micrograms below 10g**) and an SD of  $s = 6.5$ .



Evidently it's plausible in answering these questions to assume **symmetry** of the “underlying distribution”  $F$  in de Finetti's Theorem.

One standard choice, for instance, is the **Gaussian:**

$$\begin{aligned} (\mu, \sigma^2) &\sim p(\mu, \sigma^2) \\ (Y_i | \mu, \sigma^2) &\stackrel{\text{IID}}{\sim} N(\mu, \sigma^2). \end{aligned} \quad (79)$$

Here  $N(\mu, \sigma^2)$  is the familiar **normal density**

$$p(y_i | \mu, \sigma^2) = \frac{1}{\sigma\sqrt{2\pi}} \exp\left[-\frac{1}{2} \left(\frac{y_i - \mu}{\sigma}\right)^2\right]. \quad (80)$$

# Gaussian Modeling

Even though you can see from the previous graph that (79) is **not a good model** for the NB10 data, I'm going to fit it to the data for practice in working with the normal distribution from a Bayesian point of view (later we'll **improve** upon the Gaussian).

(79) is more **complicated** than the models in the AMI and LOS case studies because the parameter  $\theta$  here is a **vector**:  
 $\theta = (\mu, \sigma^2)$ .

To warm up for this new complexity let's first consider a **cut-down version of the model** in which we pretend that  $\sigma$  is known to be  $\sigma_0 = 6.5$  (the sample SD).

This **simpler model** is then

$$\left\{ \begin{array}{l} \mu \\ (Y_i|\mu) \end{array} \begin{array}{l} \sim \\ \text{IID} \\ \sim \end{array} \begin{array}{l} p(\mu) \\ N(\mu, \sigma_0^2) \end{array} \right\}. \quad (81)$$

The **likelihood function** in this model is

$$\begin{aligned} l(\mu|y) &= \prod_{i=1}^n \frac{1}{\sigma_0 \sqrt{2\pi}} \exp\left[-\frac{1}{2\sigma_0^2}(y_i - \mu)^2\right] \\ &= c \exp\left[-\frac{1}{2\sigma_0^2} \sum_{i=1}^n (y_i - \mu)^2\right] \\ &= c \exp\left[-\frac{1}{2\sigma_0^2} \left(\sum_{i=1}^n y_i^2 - 2\mu \sum_{i=1}^n y_i + n\mu^2\right)\right] \\ &= c \exp\left[-\frac{1}{2\left(\frac{\sigma_0^2}{n}\right)} (\mu - \bar{y})^2\right]. \end{aligned} \quad (82)$$

Thus the likelihood function, when thought of as a **density** for  $\mu$ , is a **normal distribution** with mean  $\bar{y}$  and SD  $\frac{\sigma_0}{\sqrt{n}}$ .

## Gaussian Modeling (continued)

Notice that this SD is the same as the frequentist **standard error** for  $\bar{Y}$  based on an IID sample of size  $n$  from the  $N(\mu, \sigma_0^2)$  distribution.

(82) also shows that the sample mean  $\bar{y}$  is a **sufficient statistic** for  $\mu$  in model (81).

In finding the conjugate prior for  $\mu$  it would be nice if the **product of two normal distributions is another normal distribution**, because that would demonstrate that the conjugate prior is normal.

Suppose therefore, to see where it leads, that the **prior for  $\mu$**  is (say)  $p(\mu) = N(\mu_0, \sigma_\mu^2)$ .

Then **Bayes' Theorem** would give

$$\begin{aligned}
 p(\mu|y) &= c p(\mu) l(\mu|y) & (83) \\
 &= c \exp\left[-\frac{1}{2\sigma_\mu^2}(\mu - \mu_0)^2\right] \exp\left[-\frac{n}{2\sigma_0^2}(\mu - \bar{y})^2\right] \\
 &= c \exp\left\{-\frac{1}{2}\left[\frac{(\mu - \mu_0)^2}{\sigma_\mu^2} + \frac{n(\mu - \bar{y})^2}{\sigma_0^2}\right]\right\},
 \end{aligned}$$

and we want this to **be of the form**

$$\begin{aligned}
 p(\mu|y) &= c \exp\left\{-\frac{1}{2}[A(\mu - B)^2 + C]\right\} \\
 &= c \exp\left\{-\frac{1}{2}[A\mu^2 - 2AB\mu + (AB^2 + C)]\right\} & (84)
 \end{aligned}$$

for some  $B, C$ , and  $A > 0$ .

Maple can help **see if this works**:

```
> collect( ( mu - mu0 )^2 / sigmamu^2 +
           n * ( mu - ybar )^2 / sigma0^2, mu );
```

$$\frac{1}{\text{sigmamu}^2} + \frac{n}{\text{sigma0}^2} \frac{(\mu - \mu_0)^2}{\mu} + \frac{-2}{\text{sigmamu}} \frac{\mu_0}{\mu} - 2 \frac{n \text{ ybar}}{\text{sigma0}^2} \frac{\mu_0}{\mu} + \frac{\mu_0^2}{\text{sigmamu}^2} + \frac{n \text{ ybar}^2}{\text{sigma0}^2}$$

# Gaussian Modeling

**Matching coefficients** for  $A$  and  $B$   
(we don't really care about  $C$ ) gives

$$A = \frac{1}{\sigma_\mu^2} + \frac{n}{\sigma_0^2} \quad \text{and} \quad B = \frac{\frac{\mu_0}{\sigma_\mu^2} + \frac{n\bar{y}}{\sigma_0^2}}{\frac{1}{\sigma_\mu^2} + \frac{n}{\sigma_0^2}}. \quad (85)$$

Since  $A > 0$  this demonstrates two things: (1) the **conjugate prior** for  $\mu$  in model (81) is **normal**, and (2) the **conjugate updating rule** (when  $\sigma_0$  is assumed known) is

$$\left\{ \begin{array}{l} \mu \sim N(\mu_0, \sigma_\mu^2) \\ (Y_i|\mu) \stackrel{\text{IID}}{\sim} N(\mu, \sigma_0^2), \\ i = 1, \dots, n \end{array} \right\} \rightarrow (\mu|y) = (\mu|\bar{y}) = N(\mu_*, \sigma_*^2), \quad (86)$$

where the **posterior mean and variance** are given by

$$\mu_* = B = \frac{\left(\frac{1}{\sigma_\mu^2}\right) \mu_0 + \left(\frac{n}{\sigma_0^2}\right) \bar{y}}{\frac{1}{\sigma_\mu^2} + \frac{n}{\sigma_0^2}} \quad \text{and} \quad \sigma_*^2 = A^{-1} = \frac{1}{\frac{1}{\sigma_\mu^2} + \frac{n}{\sigma_0^2}}. \quad (87)$$

It becomes useful in understanding the meaning of these expressions to define the **precision** of a distribution, which is just the **reciprocal** of its variance: whereas the variance and SD scales measure **uncertainty**, the precision scale quantifies **information** about an unknown.

With this convention (87) has a series of **intuitive interpretations**, as follows:

- The **prior**, considered as an **information source**, is Gaussian with mean  $\mu_0$ , variance  $\sigma_\mu^2$ , and **precision**  $\frac{1}{\sigma_\mu^2}$ , and when viewed as a data set consists of  $n_0$  (to be determined below) observations;
- The **likelihood**, considered as an **information source**, is Gaussian with mean  $\bar{y}$ , variance  $\frac{\sigma_0^2}{n}$ , and **precision**  $\frac{n}{\sigma_0^2}$ , and when viewed as a data set consists of  $n$  observations;



## Gaussian Modeling (continued)

- The **posterior**, considered as an **information source**, is Gaussian, and the posterior mean is a **weighted average** of the prior mean and data mean, with weights given by the **prior and data precisions**;
- The **posterior precision** (the reciprocal of the posterior variance) is just the **sum** of the prior and data precisions (this is why people invented the idea of precision—on this scale **knowledge** about  $\mu$  in model (81) is **additive**); and

- **Rewriting  $\mu_*$  as**

$$\mu_* = \frac{\left(\frac{1}{\sigma_\mu^2}\right) \mu_0 + \left(\frac{n}{\sigma_0^2}\right) \bar{y}}{\frac{1}{\sigma_\mu^2} + \frac{n}{\sigma_0^2}} = \frac{\left(\frac{\sigma_0^2}{\sigma_\mu^2}\right) \mu_0 + n\bar{y}}{\frac{\sigma_0^2}{\sigma_\mu^2} + n}, \quad (88)$$

you can see that the **prior sample size** is

$$n_0 = \frac{\sigma_0^2}{\sigma_\mu^2} = \frac{1}{\left(\frac{\sigma_\mu}{\sigma_0}\right)^2}, \quad (89)$$

which makes sense: the **bigger**  $\sigma_\mu$  is in relation to  $\sigma_0$ , the **less prior information** is being incorporated in the conjugate updating (86).

**Bayesian inference with multivariate  $\theta$ .** Returning now to (79) with  $\sigma^2$  unknown, (as mentioned above) this model has a ( $p = 2$ )-dimensional **parameter vector**  $\theta = (\mu, \sigma^2)$ .

When  $p > 1$  you can still use Bayes' Theorem directly to obtain the **joint posterior distribution**,

$$\begin{aligned} p(\theta|y) &= p(\mu, \sigma^2|y) = c p(\theta) l(\theta|y) \\ &= c p(\mu, \sigma^2) l(\mu, \sigma^2|y), \end{aligned} \quad (90)$$

## Multivariate Unknown $\theta$

where  $y = (y_1, \dots, y_n)$ , although making this calculation directly requires a  $p$ -dimensional **integration** to evaluate the normalizing constant  $c$ ; for example, in this case

$$\begin{aligned} c &= [p(y)]^{-1} = \left( \iint p(\mu, \sigma^2, y) d\mu d\sigma^2 \right)^{-1} \\ &= \left( \iint p(\mu, \sigma^2) l(\mu, \sigma^2 | y) d\mu d\sigma^2 \right)^{-1}. \end{aligned} \quad (91)$$

Usually, however, you'll be more interested in the **marginal posterior distributions**, in this case  $p(\mu|y)$  and  $p(\sigma^2|y)$ .

Obtaining these requires  $p$  integrations, each of dimension  $(p - 1)$ , a process that people refer to as **marginalization** or **integrating out the nuisance parameters**—for example,

$$p(\mu|y) = \int_0^\infty p(\mu, \sigma^2|y) d\sigma^2. \quad (92)$$

**Predictive** distributions also involve a  $p$ -dimensional integration: for example, with  $y = (y_1, \dots, y_n)$ ,

$$\begin{aligned} p(y_{n+1}|y) &= \iint p(y_{n+1}, \mu, \sigma^2|y) d\mu d\sigma^2 \\ &= \iint p(y_{n+1}|\mu, \sigma^2) p(\mu, \sigma^2|y) d\mu d\sigma^2. \end{aligned} \quad (93)$$

And, finally, if you're interested in a **function of the parameters**, you have some more hard integrations ahead of you.

For instance, suppose you wanted the posterior distribution for the **coefficient of variation**  $\lambda = g_1(\mu, \sigma^2) = \frac{\sqrt{\sigma^2}}{\mu}$  in model (79).

## Multivariate Unknown $\theta$

Then one fairly direct way to get this posterior (e.g., Bernardo and Smith, 1994) is to (a) introduce a **second function** of the parameters, say  $\eta = g_2(\mu, \sigma^2)$ , such that the mapping  $f = (g_1, g_2)$  from  $(\mu, \sigma^2)$  to  $(\lambda, \eta)$  is **invertible**; (b) compute the joint posterior for  $(\lambda, \eta)$  through the usual **change-of-variables formula**

$$p(\lambda, \eta|y) = p_{\mu, \sigma^2}[f^{-1}(\lambda, \eta)|y] |J_{f^{-1}}(\lambda, \eta)|, \quad (94)$$

where  $p_{\mu, \sigma^2}(\cdot, \cdot|y)$  is the joint posterior for  $\mu$  and  $\sigma^2$  and  $|J_{f^{-1}}|$  is the **determinant** of the **Jacobian** of the inverse transformation; and (c) **marginalize** in  $\lambda$  by integrating out  $\eta$  in  $p(\lambda, \eta|y)$ , in a manner analogous to (92).

Here, for instance,  $\eta = g_2(\mu, \sigma^2) = \mu$  would create an invertible  $f$ , with **inverse** defined by  $(\mu = \eta, \sigma^2 = \lambda^2 \eta^2)$ ; the **Jacobian determinant** comes out  $2\lambda\eta^2$  and (94) becomes

$$p(\lambda, \eta|y) = 2\lambda\eta^2 p_{\mu, \sigma^2}(\eta, \lambda^2\eta^2|y).$$

This process involves **two integrations**, one (of dimension  $p$ ) to get the normalizing constant that defines (94) and one (of dimension  $(p - 1)$ ) to get rid of  $\eta$ .

You can see that when  $p$  is a lot bigger than 2 all these integrals may create **severe computational problems**—this has been the **big stumbling block** for applied Bayesian work for a long time.

More than 200 years ago **Laplace** (1774)—perhaps the second applied Bayesian in history (after Bayes himself)—developed, as one avenue of solution to this problem, what people now call **Laplace approximations** to high-dimensional integrals of the type arising in Bayesian calculations (see, e.g., Tierney and Kadane, 1986).

Starting in the next case study after this one, we'll use another, computationally intensive, **simulation-based** approach: **Markov chain Monte Carlo** (MCMC).

## Gaussian Modeling

**Back to model (79).** The conjugate prior for  $\theta = (\mu, \sigma^2)$  in this model (e.g., Gelman et al., 2003) turns out to be most simply described **hierarchically**:

$$\begin{aligned}\sigma^2 &\sim \text{SI-}\chi^2(\nu_0, \sigma_0^2) \\ (\mu|\sigma^2) &\sim N\left(\mu_0, \frac{\sigma^2}{\kappa_0}\right).\end{aligned}\quad (95)$$

Here saying that  $\sigma^2 \sim \text{SI-}\chi^2(\nu_0, \sigma_0^2)$ , where SI stands for **scaled inverse**, amounts to saying that the precision  $\tau = \frac{1}{\sigma^2}$  follows a **scaled**  $\chi^2$  distribution with parameters  $\nu_0$  and  $\sigma_0^2$ .

The scaling is chosen so that  $\sigma_0^2$  can be interpreted as a **prior estimate** of  $\sigma^2$ , with  $\nu_0$  the **prior sample size** of this estimate (i.e., **think of a prior data set with  $\nu_0$  observations and sample SD  $\sigma_0$** ).

Since  $\chi^2$  is a special case of the Gamma distribution, **SI- $\chi^2$**  must be a special case of the **inverse Gamma** family—its **density** (see Gelman et al., 2003, Appendix A) is

$$\begin{aligned}\sigma^2 &\sim \text{SI-}\chi^2(\nu_0, \sigma_0^2) \leftrightarrow \\ p(\sigma^2) &= \frac{\left(\frac{1}{2}\nu_0\right)^{\frac{1}{2}\nu_0}}{\Gamma\left(\frac{1}{2}\nu_0\right)} (\sigma_0^2)^{\frac{1}{2}\nu_0} (\sigma^2)^{-(1+\frac{1}{2}\nu_0)} \exp\left(\frac{-\nu_0 \sigma_0^2}{2\sigma^2}\right).\end{aligned}\quad (96)$$

As may be verified with Maple, this distribution has **mean** (provided that  $\nu_0 > 2$ ) and **variance** (provided that  $\nu_0 > 4$ ) given by

$$E(\sigma^2) = \frac{\nu_0}{\nu_0 - 2} \sigma_0^2 \quad \text{and} \quad V(\sigma^2) = \frac{2\nu_0^2}{(\nu_0 - 2)^2(\nu_0 - 4)} \sigma_0^4. \quad (97)$$

## Gaussian Modeling (continued)

The parameters  $\mu_0$  and  $\kappa_0$  in the second level of the prior model (95),  $(\mu|\sigma^2) \sim N\left(\mu_0, \frac{\sigma^2}{\kappa_0}\right)$ , have **simple parallel interpretations** to those of  $\sigma_0^2$  and  $\nu_0$ :  $\mu_0$  is the **prior estimate** of  $\mu$ , and  $\kappa_0$  is the **prior effective sample size** of this estimate.

The **likelihood function** in model (79), with **both**  $\mu$  and  $\sigma^2$  **unknown**, is

$$\begin{aligned}
 l(\mu, \sigma^2|y) &= c \prod_{i=1}^n \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left[-\frac{1}{2\sigma^2}(y_i - \mu)^2\right] \\
 &= c (\sigma^2)^{-\frac{1}{2}n} \exp\left[-\frac{1}{2\sigma^2} \sum_{i=1}^n (y_i - \mu)^2\right] \quad (98) \\
 &= c (\sigma^2)^{-\frac{1}{2}n} \exp\left[-\frac{1}{2\sigma^2} \left(\sum_{i=1}^n y_i^2 - 2\mu \sum_{i=1}^n y_i + n\mu^2\right)\right].
 \end{aligned}$$

The **expression in brackets** in the last line of (98) is

$$\begin{aligned}
 [ \cdot ] &= -\frac{1}{2\sigma^2} \left[ \sum_{i=1}^n y_i^2 + n(\mu - \bar{y})^2 - n\bar{y}^2 \right] \quad (99) \\
 &= -\frac{1}{2\sigma^2} [n(\mu - \bar{y})^2 + (n-1)s^2],
 \end{aligned}$$

where  $s^2 = \frac{1}{n-1} \sum_{i=1}^n (y_i - \bar{y})^2$  is the **sample variance**. Thus

$$l(\mu, \sigma^2|y) = c (\sigma^2)^{-\frac{1}{2}n} \exp\left\{-\frac{1}{2\sigma^2} [n(\mu - \bar{y})^2 + (n-1)s^2]\right\},$$

and it's clear that the **vector**  $(\bar{y}, s^2)$  is **sufficient** for  $\theta = (\mu, \sigma^2)$  in this model, i.e.,  $l(\mu, \sigma^2|y) = l(\mu, \sigma^2|\bar{y}, s^2)$ .

# Gaussian Analysis

Maple can be used to make **3D** and **contour plots** of this likelihood function with the NB10 data:

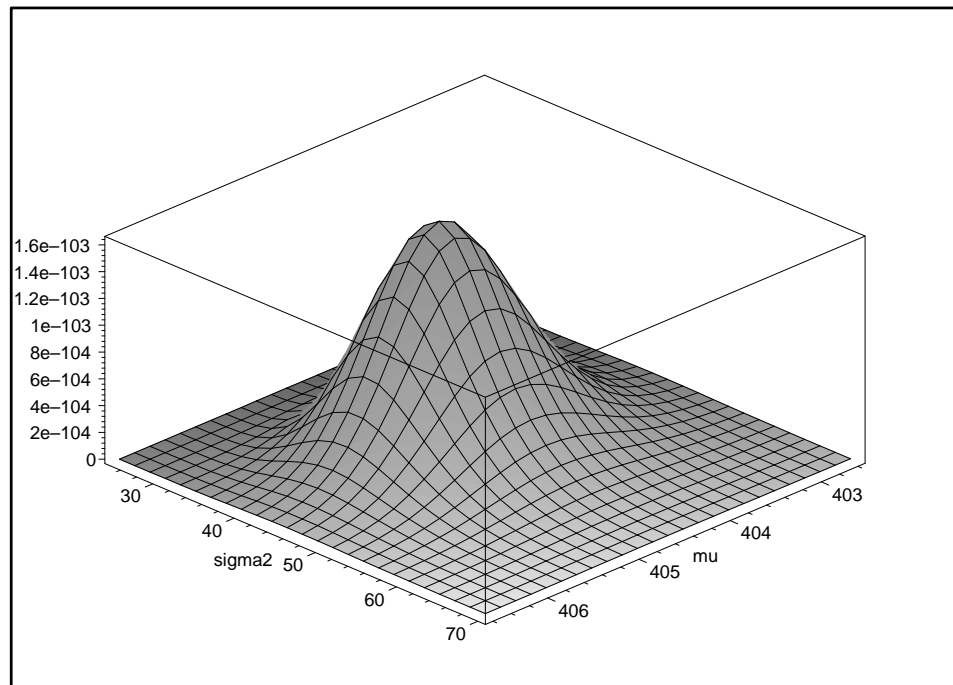
```
> l := ( mu, sigma2, ybar, s2, n ) -> sigma2^( - n / 2 ) *
    exp( - ( n * ( mu - ybar )^2 + ( n - 1 ) * s2 ) / ( 2 * sigma2 ) );
```

```
l := (mu, sigma2, ybar, s2, n) ->
```

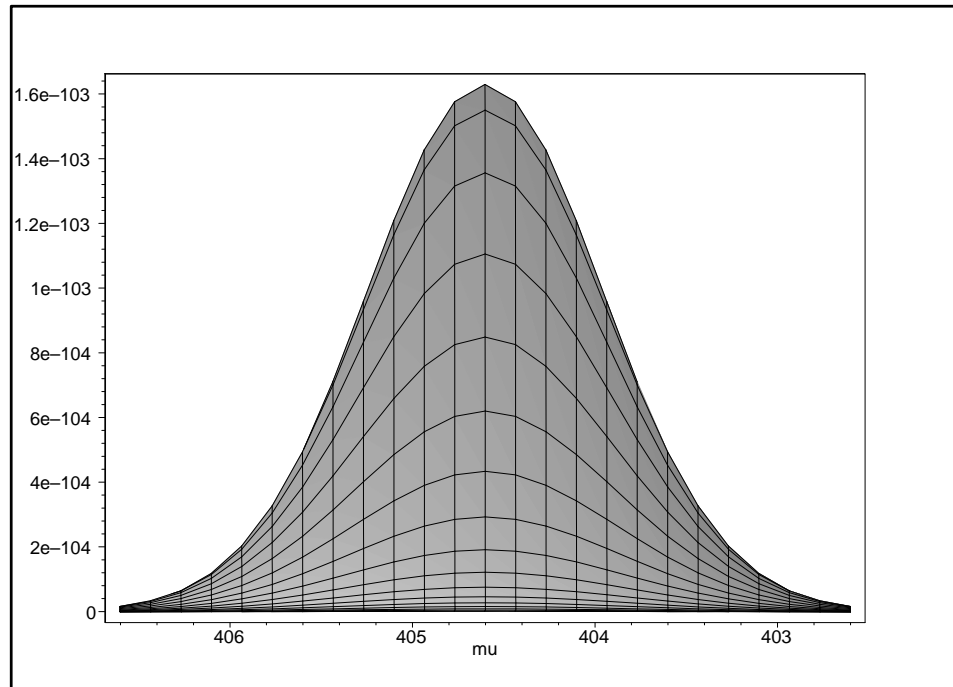
$$\text{sigma2}^{(-1/2 n)} \exp\left(-\frac{1}{2} \frac{n (\mu - \text{ybar})^2 + (n - 1) s2}{\text{sigma2}}\right)$$

```
> plotsetup( x11 );
```

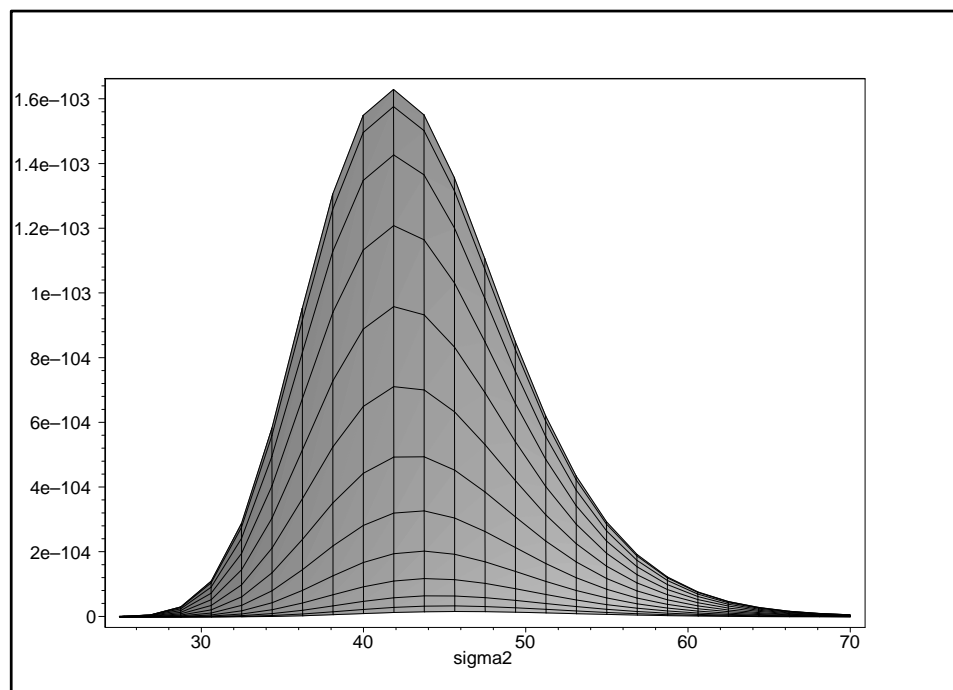
```
> plot3d( l( mu, sigma2, 404.6, 42.25, 100 ), mu = 402.6 .. 406.6,
    sigma2 = 25 .. 70 );
```



You can use the mouse to **rotate** 3D plots and get **other useful views** of them:



The **projection** or **shadow plot** of  $\mu$  looks a lot like a **normal** (or maybe a *t*) distribution.



And the shadow plot of  $\sigma^2$  looks a lot like a **Gamma** (or maybe an **inverse Gamma**) distribution.