# MAST30027: Modern Applied Statistics

# Assignment 4 Solution 2023

1. (a) Solution: The posterior distribution of the precision  $\tau$  is

$$f(\tau|x_1, \dots, x_{100}) \propto \prod_{i=1}^{100} P(x_i|\tau) f(\tau)$$

$$\propto \tau^{100/2} \exp\left\{-\frac{\tau}{2} \sum_{i=1}^{100} (x_i - 75)^2\right\} \tau^{2-1} \exp(-\tau)$$

$$\propto \tau^{52-1} \exp\left\{-\left[\frac{1}{2} \sum_{i=1}^{100} (x_i - 75)^2 + 1\right] \tau\right\}$$

This is a kernel for Gamma $(52, \frac{1}{2} \sum_{i=1}^{100} (x_i - 75)^2 + 1)$ .

Let's compute the rate parameter from the data.

> rate = 
$$sum((x-75)^2)/2 + 1$$

> rate

[1] 1805.65

Thus,  $\tau | x_1, \dots, x_{100} \sim \text{Gamma}(52, 1805.65)$ .

(b) **Solution:** Let  $\alpha^* = 52$  and  $\beta^* = 1805.65$ .

$$p(\tilde{x}|x_1, ..., x_{100}) \propto \int p(\tilde{x}|\tau) p(\tau|x_1, ..., x_{100}) d\tau$$

$$\propto \int \tau^{\frac{1}{2}} \exp\left\{-\frac{\tau}{2} (\tilde{x} - 75)^2\right\} \tau^{\alpha^* - 1} \exp(-\beta^* \tau) d\tau$$

$$\propto \int \tau^{\alpha^* + \frac{1}{2} - 1} \exp\left\{-\left[\beta^* + \frac{1}{2} (\tilde{x} - 75)^2\right] \tau\right\} d\tau$$

the integrand is proportional to the pdf of Gamma $(\alpha^* + \frac{1}{2}, \beta^* + \frac{1}{2}(\tilde{x} - 75)^2)$ 

$$\propto \frac{1}{\left(\beta^* + \frac{1}{2}(\tilde{x} - 75)^2\right)^{\alpha^* + \frac{1}{2}}}$$

$$\propto \left[1 + \frac{(\tilde{x} - 75)^2}{2\beta^*}\right]^{-(\alpha^* + \frac{1}{2})}$$

$$\propto \left[1 + \frac{(\tilde{x} - 75)^2}{2\beta^*}\right]^{-\frac{2\alpha^* + 1}{2}}.$$

Therefore we conclude that

$$\tilde{x}|x_1,...,x_{100} \sim t(\nu = 2\alpha^*, a = 75, b = \frac{\beta^*}{\alpha^*}) = t(\nu = 104, a = 75, b = 34.72404).$$

2. Posterior inference using Gibbs sampling

(a) **Solution** Let n = 100 and m = 150, we have

$$p(\mu_{1}|\mu_{2}, x_{1}, \cdots, x_{n}, y_{1}, \cdots, y_{m})$$

$$\propto p(x_{1}, \cdots, x_{n}|\mu_{1})p(y_{1}, \cdots, y_{m}|\mu_{2})p(\mu_{1}, \mu_{2})$$

$$\propto \exp\left[-\frac{1}{2}\left(\sum_{i=1}^{n}(x_{i}-\mu_{1})^{2}+2\sum_{i=1}^{m}(y_{i}-\mu_{2})^{2}+(3\mu_{1}^{2}+4\mu_{1}\mu_{2}+3\mu_{2}^{2})\right)\right]$$

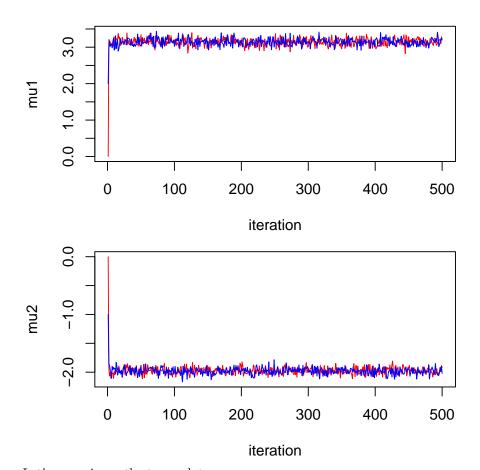
$$\propto \exp\left[-\frac{1}{2}\left((n+3)\mu_{1}^{2}-2\left(\sum_{i=1}^{n}x_{i}-2\mu_{2}\right)\mu_{1}\right)\right]$$

$$\propto \exp\left[-\frac{n+3}{2}\left(\mu_{1}^{2}-2\frac{\sum_{i=1}^{n}x_{i}-2\mu_{2}}{n+3}\mu_{1}\right)\right],$$

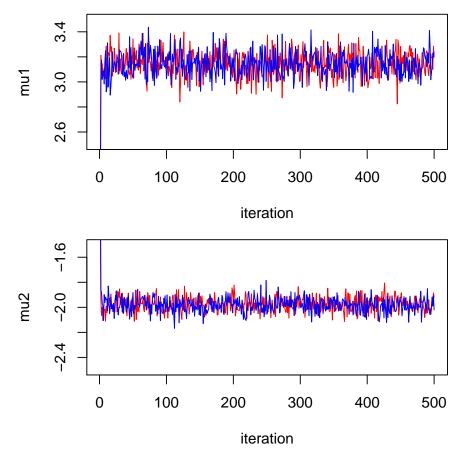
Thus, we have  $\mu_1|\mu_2, x_1, \dots, x_n, y_1, \dots, y_m \sim \text{Normal}(\frac{\sum_{i=1}^n x_i - 2\mu_2}{n+3}, \frac{1}{n+3})$ . Similarly, we have  $\mu_2|\mu_1, x_1, \dots, x_n, y_1, \dots, y_m \sim \text{Normal}(\frac{2\sum_{i=1}^m y_i - 2\mu_1}{2m+3}, \frac{1}{2m+3})$ .

(b) Solution

```
> x = scan(file="assignment4_prob2_x_2023.txt", what=double())
> y = scan(file="assignment4_prob2_y_2023.txt", what=double())
> set.seed(30027)
> # Implement Gibbs Sampler
> GibbsS <- function(mu1.0, mu2.0, nreps, x, y){
    Gsamples <- matrix(nrow=nreps, ncol=2)</pre>
    Gsamples[1,] \leftarrow c(mu1.0, mu2.0)
  # main loop
  n = length(x)
   m = length(y)
   for (i in 2:nreps) {
    mu1 = Gsamples[i-1,1]
    mu2 = Gsamples[i-1,2]
     mu1 = rnorm(1, (sum(x)-2*mu2)/(n+3), sqrt(1/(n+3)))
     mu2 = rnorm(1, (2*sum(y)-2*mu1)/(2*m+3), sqrt(1/(2*m+3)))
      Gsamples[i,] \leftarrow c(mu1, mu2)
    return(Gsamples=Gsamples)
+ }
> # number of iterations
> nreps <- 500
> # run two Gibbs sampling chains
> GibbsS1 = GibbsS(mu1.0=0, mu2.0=0, nreps, x, y)
> GibbsS2 = GibbsS(mu1.0=2, mu2.0=-1, nreps, x, y)
Make a trace plot for each of parameters.
> par(mfrow=c(2,1), mar=c(4,4,1,1))
> plot(1:nreps, GibbsS1[,1], type="l", col="red",
       ylim = c(min(GibbsS1[,1],GibbsS2[,1]), max(GibbsS1[,1],GibbsS2[,1])),
       xlab = "iteration", ylab ="mu1")
> points(1:nreps, GibbsS2[,1], type="l", col="blue")
> plot(1:nreps, GibbsS1[,2], type="l", col="red",
       ylim = c(min(GibbsS1[,2],GibbsS2[,2]), max(GibbsS1[,2],GibbsS2[,2])),
       xlab = "iteration", ylab ="mu2")
> points(1:nreps, GibbsS2[,2], type="1", col="blue")
```

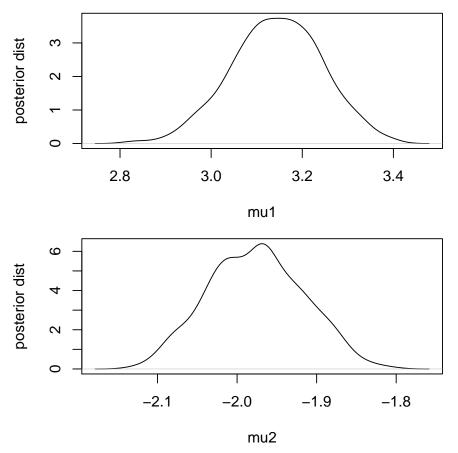


Let's zoom in on the trace plot.



The trace plots show that samples from different chains are mixed well and behave similarly.

- (c) **Solution** We will remove the first 50 samples as burn-in period.
  - 1) make a plot that shows empirical (estimated) marginal posterior distribution.
  - > par(mfrow=c(2,1), mar=c(4,4,1,1))
  - > plot(density(GibbsS1[-(1:50),1]), ylab="posterior dist", xlab="mu1", main="")
  - > plot(density(GibbsS1[-(1:50),2]), ylab="posterior dist", xlab="mu2", main="")



2) estimate marginal posterior mean.

```
> # for mu1
> mean(GibbsS1[-(1:50),1])
[1] 3.144036
> # for mu2
```

> mean(GibbsS1[-(1:50),2])

[1] -1.976291

3) report a 90% credible interval for the marginal posterior distribution.

> quantile(GibbsS1[-(1:50),1], probs=c(0.05, 0.95))

```
5 quantile(GibbsSi[-(1:50),1], probs-c(0.05, 0.95))
5% 95%
2.974518 3.306848
> # for mu2
> quantile(GibbsSi[-(1:50),2], probs=c(0.05, 0.95))
5% 95%
-2.076106 -1.876740
```

## Posterior inference using the Metropolis-Hastings (MH) algorithm

(d) **Solution** The posterior distribution is proportional to the product of the likelihood and prior distribution. Thus, we will use the product of the likelihood and prior distribution as a target distribution  $\pi(\mu_1, \mu_2)$  in the Metropolis-Hastings algorithm.

```
> # log likelihood
> likelihood <- function(param, x, y){</pre>
```

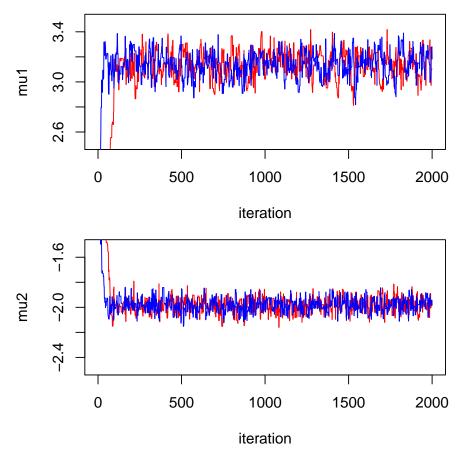
```
mu1 = param[1]
+
      mu2 = param[2]
      singlelikelihoodsX = dnorm(x, mean = mu1, sd = 1, log = T)
      singlelikelihoodsY = dnorm(y, mean = mu2, sd = sqrt(1/2), log = T)
      sumll = sum(singlelikelihoodsX) + sum(singlelikelihoodsY)
+
      return(sumll)
+ }
> # log prior
> prior <- function(param){</pre>
     mu1 = param[1]
     mu2 = param[2]
+
      return(-0.5*(3*mu1^2 + 4*mu1*mu2 + 3*mu2^2))
+ }
> # log posterior distribution up to a constant
> posterior <- function(param, x, y){</pre>
     return (likelihood(param, x, y) + prior(param))
+ }
> # proposal function
> proposalfunction <- function(param, delta){</pre>
  mu1 = param[1]
  mu2 = param[2]
   new.mu1 = rnorm(1, mu1, delta)
  new.mu2 = rnorm(1, mu2, delta)
+
   return(as.vector(c(new.mu1, new.mu2)))
+ }
> # log prob of proposal function
> proposal.prob <- function(old.param, new.param, delta){</pre>
   mu1 = old.param[1]
   mu2 = old.param[2]
+
   new.mu1 = new.param[1]
   new.mu2 = new.param[2]
   return(dnorm(new.mu1, mean = mu1, sd = delta, log = T) +
             dnorm(new.mu2, mean = mu2, sd= delta, log=T))
+ }
> # Metropolis algorithm
> run_metropolis_MCMC <- function(x, y, startvalue, iterations, delta){
      chain = array(dim = c(iterations, 2))
+
+
      chain[1,] = startvalue
      for (i in 1:(iterations-1)){
          proposal = proposalfunction(chain[i,], delta)
+
          probab = exp(posterior(proposal,x, y) +
+
                         proposal.prob(proposal, chain[i,], delta) -
                         posterior(chain[i,], x, y) -
                         proposal.prob(chain[i,], proposal, delta))
+
          if (runif(1) < probab){</pre>
              chain[i+1,] = proposal
          }else{
              chain[i+1,] = chain[i,]
          }
      }
      return(chain)
```

```
> # number of iterations
> nreps <- 2000
> # run two MH chains
> MHS1 = run_metropolis_MCMC(x, y, startvalue = c(0,0), nreps, delta = 1/10)
> MHS2 = run_metropolis_MCMC(x, y, startvalue = c(2,-1), nreps, delta = 1/10)
Make a trace plot for each of parameters.
> par(mfrow=c(2,1), mar=c(4,4,1,1))
> plot(1:nreps, MHS1[,1], type="l", col="red",
       ylim = c(min(MHS1[,1],MHS2[,1]), max(MHS1[,1],MHS2[,1])),
       xlab = "iteration", ylab ="mu1")
> points(1:nreps, MHS2[,1], type="l", col="blue")
> plot(1:nreps, MHS1[,2], type="l", col="red",
       ylim = c(min(MHS1[,2],MHS2[,2]), max(MHS1[,2],MHS2[,2])),
       xlab = "iteration", ylab ="mu2")
> points(1:nreps, MHS2[,2], type="1", col="blue")
     0
     \alpha
     1.0
            0
                       500
                                   1000
                                               1500
                                                           2000
                                 iteration
            0
                       500
                                   1000
                                               1500
                                                           2000
                                 iteration
Let's zoom in on the trace plot.
> par(mfrow=c(2,1), mar=c(4,4,1,1))
> plot(1:nreps, MHS1[,1], type="l", col="red", ylim = c(2.5, 3.5),
       xlab = "iteration", ylab ="mu1")
```

> plot(1:nreps, MHS1[,2], type="l", col="red", ylim = c(-2.5, -1.5),

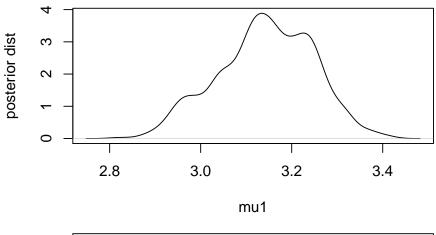
> points(1:nreps, MHS2[,1], type="l", col="blue")

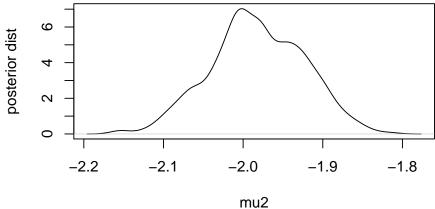
+ xlab = "iteration", ylab ="mu2")
> points(1:nreps, MHS2[,2], type="1", col="blue")



The trace plots show that samples from different chains are mixed well and behave similarly after 2000 iterations.

- (e) Solution We will remove the first 300 samples as burn-in period.
  - 1) make a plot that shows empirical (estimated) marginal posterior distribution.
  - > par(mfrow=c(2,1), mar=c(4,4,1,1))
  - > plot(density(MHS1[-(1:300),1]), ylab="posterior dist", xlab="mu1", main="")
  - > plot(density(MHS1[-(1:300),2]), ylab="posterior dist", xlab="mu2", main="")





- 2) estimate marginal posterior mean.
- > # for mu1
- > mean(MHS1[-(1:300),1])
- [1] 3.142711
- > # for mu2
- > mean(MHS1[-(1:300),2])
- [1] -1.982071
- 3) report a 90% credible interval for the marginal posterior distribution.
- > # for mu1
- > quantile(MHS1[-(1:300),1], probs=c(0.05, 0.95))

- 2.955025 3.305431
- > # for mu2
- > quantile(MHS1[-(1:300),2], probs=c(0.05, 0.95))

-2.081424 -1.889318

### Posterior inference using Variational Inference (VI)

(f) **Solution** From problem 1, we have

$$\log p(\mu_1, \mu_2, x_1, \dots, x_n, y_1, \dots, y_m)$$

$$= -\frac{1}{2} \left( \sum_{i=1}^n (x_i - \mu_1)^2 + 2 \sum_{i=1}^m (y_i - \mu_2)^2 + (3\mu_1^2 + 4\mu_1\mu_2 + 3\mu_2^2) \right) + \text{const}$$

Then

$$\log q_{\mu_1}^*(\mu_1) = -\frac{1}{2} E_{\mu_2} \left[ \sum_{i=1}^n (x_i - \mu_1)^2 + 2 \sum_{i=1}^m (y_i - \mu_2)^2 + (3\mu_1^2 + 4\mu_1\mu_2 + 3\mu_2^2) \right] + \text{const}$$

$$= -\frac{1}{2} \left( \sum_{i=1}^n (x_i - \mu_1)^2 + E_{\mu_2} \left[ 2 \sum_{i=1}^m (y_i - \mu_2)^2 \right] + 3\mu_1^2 + 4\mu_1 E_{\mu_2}(\mu_2) \right) + \text{const}$$

$$= -\frac{1}{2} \left( (n+3)\mu_1^2 - 2 \left( \sum_{i=1}^n x_i - 2E_{\mu_2}(\mu_2) \right) \mu_1 \right) + \text{const}$$

Hence,  $q_{\mu_1}^*(\mu_1)$  is the pdf of  $N(\mu_1^*, \sigma_1^{2*})$ , where  $\mu_1^* = \frac{\sum_{i=1}^n x_i - 2E_{\mu_2}(\mu_2)}{n+3}$  and  $\sigma_1^{2*} = \frac{1}{n+3}$ . Similarly,

$$\log q_{\mu_2}^*(\mu_2) = -\frac{1}{2} E_{\mu_1} \left[ \sum_{i=1}^n (x_i - \mu_1)^2 + 2 \sum_{i=1}^m (y_i - \mu_2)^2 + (3\mu_1^2 + 4\mu_1\mu_2 + 3\mu_2^2) \right] + \text{const}$$

$$= -\frac{1}{2} \left( E_{\mu_1} \left[ \sum_{i=1}^n (x_i - \mu_1)^2 \right] + 2 \sum_{i=1}^m (y_i - \mu_2)^2 + 4 E_{\mu_1}(\mu_1) \mu_2 + 3\mu_2^2 \right) + \text{const}$$

$$= -\frac{1}{2} \left( (2m+3)\mu_2^2 - 2 \left( 2 \sum_{i=1}^m y_i - 2E_{\mu_1}(\mu_1) \right) \mu_2 \right) + \text{const}$$

Hence,  $q_{\mu_2}^*(\mu_2)$  is the pdf of  $N(\mu_2^*, \sigma_2^{2*})$ , where  $\mu_2^* = \frac{2\sum_{i=1}^m y_i - 2E_{\mu_1}(\mu_1)}{2m+3}$  and  $\sigma_2^{2*} = \frac{1}{2m+3}$ .

#### (g) Solution

$$ELBO(q_{\mu_1}^*(\mu_1), q_{\mu_2}^*(\mu_2)) = ELBO(\mu_1^*, \mu_2^*)$$

$$= E_{\mu_1, \mu_2} \left[ \log p(\mu_1, \mu_2, x_1, \cdots, x_n, y_1, \cdots, y_m) \right] - E_{\mu_1, \mu_2} \left[ \log(q_{\mu_1}^*(\mu_1)) \right] - E_{\mu_1, \mu_2} \left[ \log(q_{\mu_2}^*(\mu_2)) \right]$$

Each term is computed as follow,

$$\begin{split} &E_{\mu_1,\mu_2}\left[\log p(\mu_1,\mu_2,x_1,\cdots,x_n,y_1,\cdots,y_m)\right] \\ &= -\frac{1}{2}E_{\mu_1,\mu_2}\left[\sum_{i=1}^n(x_i-\mu_1)^2+2\sum_{i=1}^m(y_i-\mu_2)^2+(3\mu_1^2+4\mu_1\mu_2+3\mu_2^2)\right] + \text{const} \\ &= -\frac{1}{2}\left[\sum_{i=1}^nE_{\mu_1}(x_i-\mu_1)^2+2\sum_{i=1}^mE_{\mu_2}(y_i-\mu_2)^2+3(\sigma_1^{2*}+\mu_1^{*2})+4\mu_1^*\mu_2^*+3(\sigma_2^{2*}+\mu_2^{*2})\right] + \text{const} \\ &= -\frac{1}{2}\left[\sum_{i=1}^n\left(\sigma_1^{2*}+\mu_1^{*2}-2x_i\mu_1^*\right)+2\sum_{i=1}^m\left(\sigma_2^{2*}+\mu_2^{*2}-2y_i\mu_2^*\right)+3(\sigma_1^{2*}+\mu_1^{*2})+4\mu_1^*\mu_2^*+3(\sigma_2^{2*}+\mu_2^{*2})\right] + \text{const} \end{split}$$

$$E_{\mu_1,\mu_2} \left[ \log(q_{\mu_1}^*(\mu_1)) \right]$$

$$= -\frac{1}{2} E_{\mu_1,\mu_2} \left[ \frac{(\mu_1 - \mu_1^*)^2}{\sigma_1^{2*}} \right] + \text{const}$$

$$= -\frac{1}{2} + \text{const}$$

$$E_{\mu_1,\mu_2} \left[ \log(q_{\mu_2}^*(\mu_2)) \right]$$

$$= -\frac{1}{2} E_{\mu_1,\mu_2} \left[ \frac{(\mu_2 - \mu_2^*)^2}{\sigma_2^{2*}} \right] + \text{const}$$

$$= -\frac{1}{2} + \text{const}$$

#### (h) **Solution**

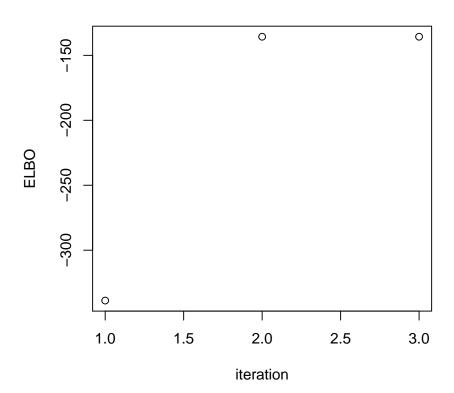
Implementation CAVI algorithm

```
> # x y : data
> # initial values for mu1*, sigma1.2*, mu2*, sigma2.2*:
             mu1.vi.init, sigma1.2.vi.init, mu2.vi.init, sigma2.2.vi.init
> # epsilon : If the ELBO has changed by less than epsilon,
             the CAVI algorithm will stop
> # max.iter: maximum number of iteration
> cavi.normal <- function(x, y, mu1.vi.init, sigma1.2.vi.init,
                          mu2.vi.init, sigma2.2.vi.init, epsilon=1e-5, max.iter=100){
+
+
   n = length(x)
   m = length(y)
   mu1.vi = mu1.vi.init
   sigma1.2.vi = sigma1.2.vi.init
   mu2.vi = mu2.vi.init
   sigma2.2.vi = sigma2.2.vi.init
    # store the ELBO for each iteration
   elbo = c()
   # I will store mu1*, sigma1.2*, mu2*, sigma2.2* for each iteration
   mu1.vi.list = sigma1.2.vi.list = mu2.vi.list = sigma2.2.vi.list = c()
    # compute the ELBO using initial values of mu1*, sigma1.2*, mu2*, sigma2.2*
   Elogq.mu1 = Elogq.mu2 = -1/2
    # Elogp.x.y.mu1.mu2
   A = sigma1.2.vi + mu1.vi^2 - 2*x*mu1.vi + x*x
   B = sigma2.2.vi + mu2.vi^2 - 2*y*mu2.vi + y*y
   Elogp.x.y.mu1.mu2 = -0.5*sum(A) - sum(B) -
      0.5*(3*(sigma1.2.vi + mu1.vi^2) + 4*mu1.vi*mu2.vi + 3*(sigma2.2.vi + mu2.vi^2))
+
    elbo = c(elbo, Elogp.x.y.mu1.mu2 -Elogq.mu1 - Elogq.mu2)
   mu1.vi.list = c(mu1.vi.list, mu1.vi)
   sigma1.2.vi.list = c(sigma1.2.vi.list, sigma1.2.vi)
   mu2.vi.list = c(mu2.vi.list, mu2.vi)
   sigma2.2.vi.list = c(sigma2.2.vi.list, sigma2.2.vi)
    # set the change in the ELBO with 1
+
   delta.elbo = 1
    # number of iteration
   n.iter = 1
+
+
    # If the elbo has changed by less than epsilon, the CAVI will stop.
    while((delta.elbo > epsilon) & (n.iter <= max.iter)){</pre>
+
      # Update mu1.vi and sigma1.2.vi
      mu1.vi = (sum(x) - 2*mu2.vi)/(n + 3)
      sigma1.2.vi = 1/(n+3)
      # Update mu2.vi and sigma2.2.vi
      mu2.vi = (2*sum(y) - 2*mu1.vi)/(2*m + 3)
```

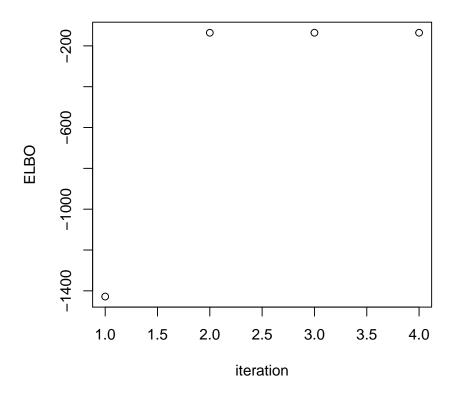
```
+
+
      # compute the ELBO using the current values of mu1*, sigma1.2*, mu2*, sigma2.2*
      Elogq.mu1 = Elogq.mu2 = -1/2
+
      # Elogp.x.y.mu1.mu2
+
      A = sigma1.2.vi + mu1.vi^2 - 2*x*mu1.vi + x*x
      B = sigma2.2.vi + mu2.vi^2 - 2*y*mu2.vi + y*y
+
      Elogp.x.y.mu1.mu2 = -0.5*sum(A) - sum(B) -
+
        0.5*(3*(sigma1.2.vi + mu1.vi^2) + 4*mu1.vi*mu2.vi + 3*(sigma2.2.vi + mu2.vi^2))
+
+
      elbo = c(elbo, Elogp.x.y.mu1.mu2 -Elogq.mu1 - Elogq.mu2)
+
      mu1.vi.list = c(mu1.vi.list, mu1.vi)
      sigma1.2.vi.list = c(sigma1.2.vi.list, sigma1.2.vi)
+
      mu2.vi.list = c(mu2.vi.list, mu2.vi)
      sigma2.2.vi.list = c(sigma2.2.vi.list, sigma2.2.vi)
+
+
      # compute the change in the elbo
      delta.elbo = elbo[length(elbo)] - elbo[length(elbo)-1]
+
      # increase the number of iteration
+
      n.iter = n.iter + 1
+
   return(list(elbo = elbo,
                mu1.vi.list = mu1.vi.list, sigma1.2.vi.list=sigma1.2.vi.list,
+
                mu2.vi.list = mu2.vi.list, sigma2.2.vi.list=sigma2.2.vi.list))
+ }
Applying the implemented algorithm. Run the CAVI algorithm with different initial
values and check that the ELBO increases at each step by plotting them.
> cavi1 = cavi.normal(x, y,
                      mu1.vi.init = 3, sigma1.2.vi.init = 1,
                      mu2.vi.init = -2, sigma2.2.vi.init = 1,
                      epsilon=1e-5, max.iter=100)
> cavi.res = cavi1
> cavi.res$elbo
[1] -338.7661 -135.6699 -135.6699
> plot(cavi.res$elbo, ylab='ELBO', xlab='iteration')
> print(paste("mu1* and sigma1.2* = (",
              round(cavi.res$mu1.vi.list[length(cavi.res$mu1.vi.list)],2), ",",
              round(cavi.res$sigma1.2.vi.list[length(cavi.res$sigma1.2.vi.list)],4),
+
              ")", sep=""))
[1] "mu1* and sigma1.2* = (3.14,0.0097)"
> print(paste("mu2* and sigma2.2* = (",
              round(cavi.res$mu2.vi.list[length(cavi.res$mu2.vi.list)],2), ",",
+
              round(cavi.res$sigma2.2.vi.list[length(cavi.res$sigma2.2.vi.list)],4),
              ")", sep=""))
[1] "mu2* and sigma2.2* = (-1.98, 0.0033)"
```

sigma2.2.vi = 1/(2\*m + 3)

+



```
> cavi2 = cavi.normal(x, y,
                      mu1.vi.init = 0, sigma1.2.vi.init = 1,
                      mu2.vi.init = 0, sigma2.2.vi.init = 1,
                      epsilon=1e-5, max.iter=100)
> cavi.res = cavi2
> cavi.res$elbo
[1] -1428.0667 -135.7461 -135.6699 -135.6699
> plot(cavi.res$elbo, ylab='ELBO', xlab='iteration')
> print(paste("mu1* and sigma1.2* = (",
              round(cavi.res$mu1.vi.list[length(cavi.res$mu1.vi.list)],2), ",",
+
              round(cavi.res$sigma1.2.vi.list[length(cavi.res$sigma1.2.vi.list)],4),
              ")", sep=""))
[1] "mu1* and sigma1.2* = (3.14, 0.0097)"
> print(paste("mu2* and sigma2.2* = (",
              round(cavi.res$mu2.vi.list[length(cavi.res$mu2.vi.list)],2), ",",
              round(cavi.res$sigma2.2.vi.list[length(cavi.res$sigma2.2.vi.list)],4),
              ")", sep=""))
[1] "mu2* and sigma2.2* = (-1.98, 0.0033)"
```



The two CAVI runs have equally highest ELBO. You can see that approximated posterior distributions from the runs are the same. I will use the output from the first run:  $q_{\mu_1}^*(\mu_1)$  is a pdf of N(3.14,0.0097) and  $q_{\mu_2}^*(\mu_2)$  is a pdf of N(-1.98,0.0033).