## Computing Homework 1

Code ▼

Due: Wednesday 11/8 at 11:59pm, Canvas submission (.nb.html file only)

#### Honor Pledge

On my honor, I have neither received nor given any unauthorized assistance on this Homework.

SIGNED: 218007361

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# Load packages
library(ggplot2)
library(gridExtra)
library(rstan)
library(bayesplot)

In this exercise we consider two methods of simulating from the posterior distribution of a Beta-Binomial Bayesian model. The prior distribution is

$$\pi \sim Beta(1,2),$$

and the data model is

$$y_i \mid \pi \stackrel{ind}{\sim} Bin(6,\pi).$$

Two observations are made:

$$ec{y}=(y_1,y_2)=(3,4).$$

Complete the following questions.

## Question 1: analytical posterior (1pt)

Fill in the two blanks with numbers: the analytical posterior for  $\pi$  is

$$\pi \mid ec{y} \sim Beta(8,7).$$

## Question 2: Posterior simulation in rstan (5 pts total)

#### 2 - 1. (1pt)

Fill in the code chunk named <code>model</code> below, to define the Bayesian model structure in <code>rstan</code>. Use the following notation in your definition:

- Y for the data vector  $\vec{y}$ ;
- pi for the parameter.

Be sure to specify the data type, range, vector length (if applicable), etc.

```
the_model <- "
  data {
    int<lower=0, upper=6> Y[2];
  }
  parameters {
    real<lower=0, upper=1> pi;
  }
  model {
    Y ~ binomial(6, pi);
    pi ~ beta(1, 2);
  }
  "
```

### 2 - 2. (0pt)

Complete the code below to simulate from the posterior using the model defined above. Use the following settings:

- number of chains: 4
- number of iter: 10,000.

```
SAMPLING FOR MODEL 'anon model' NOW (CHAIN 1).
Chain 1:
Chain 1: Gradient evaluation took 5e-06 seconds
Chain 1: 1000 transitions using 10 leapfrog steps per transition would take 0.05 seconds.
Chain 1: Adjust your expectations accordingly!
Chain 1:
Chain 1:
Chain 1: Iteration:
                       1 / 10000 [ 0%]
                                         (Warmup)
Chain 1: Iteration: 1000 / 10000 [ 10%]
                                        (Warmup)
Chain 1: Iteration: 2000 / 10000 [ 20%]
                                        (Warmup)
Chain 1: Iteration: 3000 / 10000 [ 30%] (Warmup)
Chain 1: Iteration: 4000 / 10000 [ 40%] (Warmup)
Chain 1: Iteration: 5000 / 10000 [ 50%]
                                        (Warmup)
Chain 1: Iteration: 5001 / 10000 [ 50%] (Sampling)
Chain 1: Iteration: 6000 / 10000 [ 60%] (Sampling)
Chain 1: Iteration: 7000 / 10000 [ 70%]
                                        (Sampling)
Chain 1: Iteration: 8000 / 10000 [ 80%] (Sampling)
Chain 1: Iteration: 9000 / 10000 [ 90%] (Sampling)
Chain 1: Iteration: 10000 / 10000 [100%]
                                          (Sampling)
Chain 1:
Chain 1: Elapsed Time: 0.029 seconds (Warm-up)
Chain 1:
                        0.034 seconds (Sampling)
Chain 1:
                        0.063 seconds (Total)
Chain 1:
SAMPLING FOR MODEL 'anon model' NOW (CHAIN 2).
Chain 2:
Chain 2: Gradient evaluation took 3e-06 seconds
Chain 2: 1000 transitions using 10 leapfrog steps per transition would take 0.03 seconds.
Chain 2: Adjust your expectations accordingly!
Chain 2:
Chain 2:
Chain 2: Iteration:
                       1 / 10000 [ 0%] (Warmup)
Chain 2: Iteration: 1000 / 10000 [ 10%] (Warmup)
Chain 2: Iteration: 2000 / 10000 [ 20%] (Warmup)
Chain 2: Iteration: 3000 / 10000 [ 30%]
                                        (Warmup)
Chain 2: Iteration: 4000 / 10000 [ 40%] (Warmup)
Chain 2: Iteration: 5000 / 10000 [ 50%]
                                        (Warmup)
Chain 2: Iteration: 5001 / 10000 [ 50%] (Sampling)
Chain 2: Iteration: 6000 / 10000 [ 60%] (Sampling)
Chain 2: Iteration: 7000 / 10000 [ 70%]
                                         (Sampling)
Chain 2: Iteration: 8000 / 10000 [ 80%]
                                         (Sampling)
Chain 2: Iteration: 9000 / 10000 [ 90%]
                                        (Sampling)
Chain 2: Iteration: 10000 / 10000 [100%]
                                         (Sampling)
Chain 2:
Chain 2: Elapsed Time: 0.028 seconds (Warm-up)
Chain 2:
                        0.178 seconds (Sampling)
Chain 2:
                        0.206 seconds (Total)
Chain 2:
SAMPLING FOR MODEL 'anon_model' NOW (CHAIN 3).
```

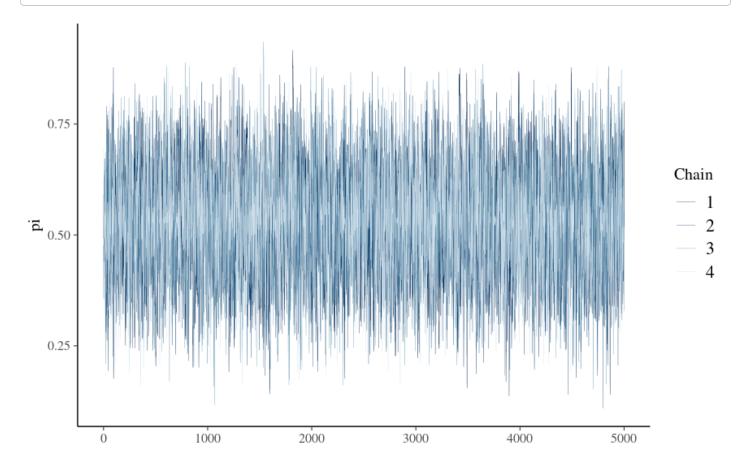
```
Chain 3:
Chain 3: Gradient evaluation took 3e-06 seconds
Chain 3: 1000 transitions using 10 leapfrog steps per transition would take 0.03 seconds.
Chain 3: Adjust your expectations accordingly!
Chain 3:
Chain 3:
Chain 3: Iteration:
                       1 / 10000 [ 0%]
                                         (Warmup)
Chain 3: Iteration: 1000 / 10000 [ 10%]
                                         (Warmup)
Chain 3: Iteration: 2000 / 10000 [ 20%] (Warmup)
Chain 3: Iteration: 3000 / 10000 [ 30%]
                                         (Warmup)
Chain 3: Iteration: 4000 / 10000 [ 40%]
                                         (Warmup)
Chain 3: Iteration: 5000 / 10000 [ 50%] (Warmup)
Chain 3: Iteration: 5001 / 10000 [ 50%]
                                         (Sampling)
Chain 3: Iteration: 6000 / 10000 [ 60%] (Sampling)
Chain 3: Iteration: 7000 / 10000 [ 70%]
                                        (Sampling)
Chain 3: Iteration: 8000 / 10000 [ 80%]
                                         (Sampling)
Chain 3: Iteration: 9000 / 10000 [ 90%] (Sampling)
Chain 3: Iteration: 10000 / 10000 [100%] (Sampling)
Chain 3:
Chain 3: Elapsed Time: 0.029 seconds (Warm-up)
Chain 3:
                        0.029 seconds (Sampling)
Chain 3:
                        0.058 seconds (Total)
Chain 3:
SAMPLING FOR MODEL 'anon_model' NOW (CHAIN 4).
Chain 4:
Chain 4: Gradient evaluation took 3e-06 seconds
Chain 4: 1000 transitions using 10 leapfrog steps per transition would take 0.03 seconds.
Chain 4: Adjust your expectations accordingly!
Chain 4:
Chain 4:
Chain 4: Iteration:
                       1 / 10000 [ 0%]
                                        (Warmup)
Chain 4: Iteration: 1000 / 10000 [ 10%]
                                         (Warmup)
Chain 4: Iteration: 2000 / 10000 [ 20%] (Warmup)
Chain 4: Iteration: 3000 / 10000 [ 30%]
                                        (Warmup)
Chain 4: Iteration: 4000 / 10000 [ 40%] (Warmup)
Chain 4: Iteration: 5000 / 10000 [ 50%] (Warmup)
Chain 4: Iteration: 5001 / 10000 [ 50%]
                                         (Sampling)
Chain 4: Iteration: 6000 / 10000 [ 60%] (Sampling)
Chain 4: Iteration: 7000 / 10000 [ 70%]
                                        (Sampling)
Chain 4: Iteration: 8000 / 10000 [ 80%]
                                         (Sampling)
Chain 4: Iteration: 9000 / 10000 [ 90%]
                                         (Sampling)
Chain 4: Iteration: 10000 / 10000 [100%]
                                         (Sampling)
Chain 4:
Chain 4: Elapsed Time: 0.029 seconds (Warm-up)
Chain 4:
                        0.03 seconds (Sampling)
Chain 4:
                        0.059 seconds (Total)
Chain 4:
```

#### 2 - 3. (1pt)

Visualize the trace plots of the chains in the\_sim:

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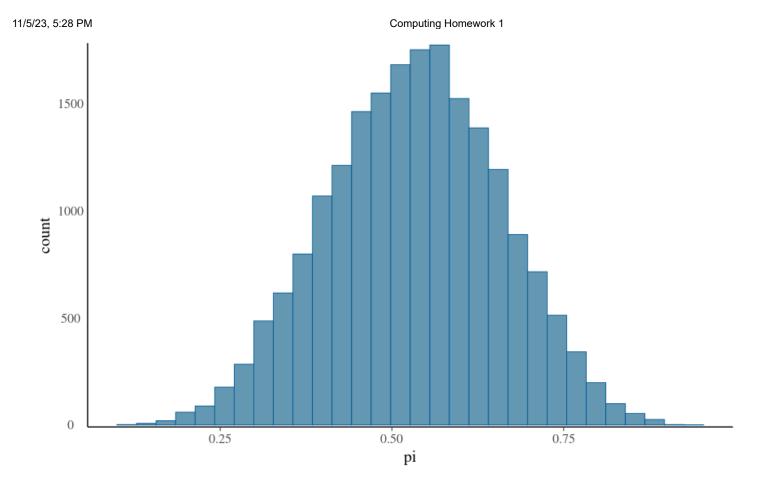
```
mcmc_trace(the_sim, pars = "pi", size = 0.1)
```



## 2 - 4. (1pt)

Visualize the histogram of the chains in the\_sim:

```
mcmc_hist(the_sim, pars = "pi") +
  yaxis_text(TRUE) + ylab("count")
```



## 2 - 5. (1pt)

Calculate the effective sample size ratio for the\_sim:

neff\_ratio(the\_sim, pars = c("pi"))

[1] 0.3701599

· Is this result satisfactory? Yes

## 2 - 6. (1pt)

Calculate the  $\hat{R}$  for the chains in <code>the\_sim</code>:

rhat(the\_sim, pars = "pi")

[1] 1.000401

· Is this result satisfactory? Yes

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# Question 3: Posterior simulation via Independence Sampler (4 pts total)

Now forget about what you did using rstan. Instead, use the Independence Sampler algorithm to target the posterior distribution for  $\pi$ , assuming you do not know its analytical distribution. Use proposal distribution

$$\pi^* \sim Beta(a,b).$$

#### 3 - 1. (1pt)

Complete the code chunk below to define the function, <code>one\_iteration</code>, that produces one iteration of the independent sampler. Here,

- a : hyperparameter a;
- b : hyperparameter b;
- current : current value of the chain.

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#### 3 - 2. (3 pts)

The function whole\_tour, which produces N iterations of the independence sampler, has been defined for you below.

```
whole_tour <- function(N, a, b){</pre>
 # 1. Start the chain at location 0.5
 current <- 0.5
 # 2. Initialize the simulation
 pi <- rep(0, N)
 # 3. Simulate N Markov chain stops
 for(i in 1:N){
    # Simulate one iteration
    sim <- one_iteration(a = a, b = b, current = current)</pre>
    # Record next location
    pi[i] <- sim$next_stop</pre>
    # Reset the current location
    current <- sim$next_stop</pre>
 # 4. Return the chain locations
  return(data.frame(iteration = c(1:N), pi))
}
```

Complete the following chunks of code to:

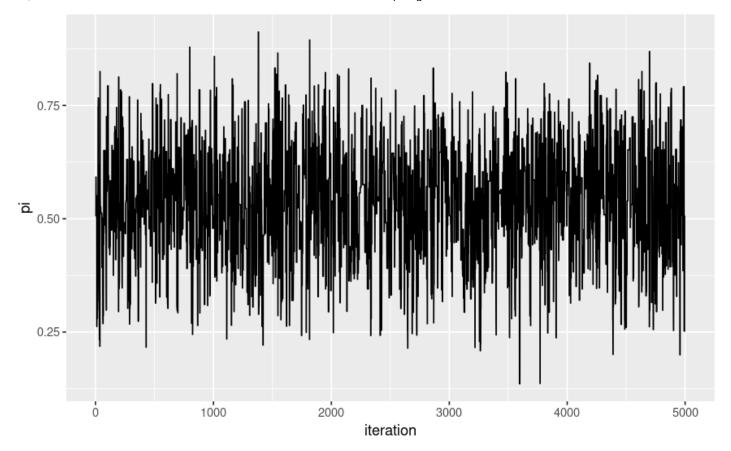
• (1pt) Use the whole\_tour function to produce the object independence\_sim, consisting of N = 5000 iterations from the independent sampler, with prior hyperparameters a = 1 and b = 2;

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```
set.seed(1000)
independence_sim <- whole_tour(N=5000, a=1, b=2)</pre>
```

• (1pt) Produce the trace plot of the chain in independence\_sim:

```
ggplot(independence_sim, aes(x = iteration, y = pi)) +
  geom_line()
```



• (1pt) Produce the histogram of the chain in independence\_sim, and overlay it with the analytical posterior for  $\pi$  which you derived from Question 1.

```
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```

```
ggplot(independence_sim, aes(x = pi)) +
  geom_histogram(aes(y = ..density..), bins=30, color = "white") +
  stat_function(fun = dbeta, args = list(8,7), color = "blue")
```

