

Bayesian HW5

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11/10/2019

BDA Problem 11.2

Metropolis algorithm: Replicate the computations for the bioassay example of Section 3.7 using the Metropolis algorithm. Be sure to define your starting points and your jumping rule. Compute with log-densities (see page 261). Run the simulations long enough for approximate convergence.

```
a.c = 0.8
b.c = 7.7
a.sd = 1
b.sd = 4.9

mcmc_array <- function (ns, nchains = 1, params) {
  nparams <- length(params)
  array(dim = c(ns, nchains, nparams),
        dimnames = list(iterations = NULL,
                          chains = paste0("chain", 1:nchains),
                          parameters = params))
}

data = data.frame("y" = c(0,1,3,5),
                  "n" = c(5,5,5,5),
                  "x" = c(-0.86,-0.3,-0.05,0.73))

in.lo = function(al,be){
  the = al + be * data$x
  the = invlogit(the)
  return(the)
}

po = function(al,be){
  the = in.lo(al=al,be=be)
  sum.cho = sum(lchoose(n = data$n,k = data$y))
  p2 = sum(data$y*log(the))
  p3 = sum(log(1-the)*(data$n-data$y))
  lopo = exp(sum.cho + p2 + p3)
  return(lopo)
}

nc = 4
ns = 10000

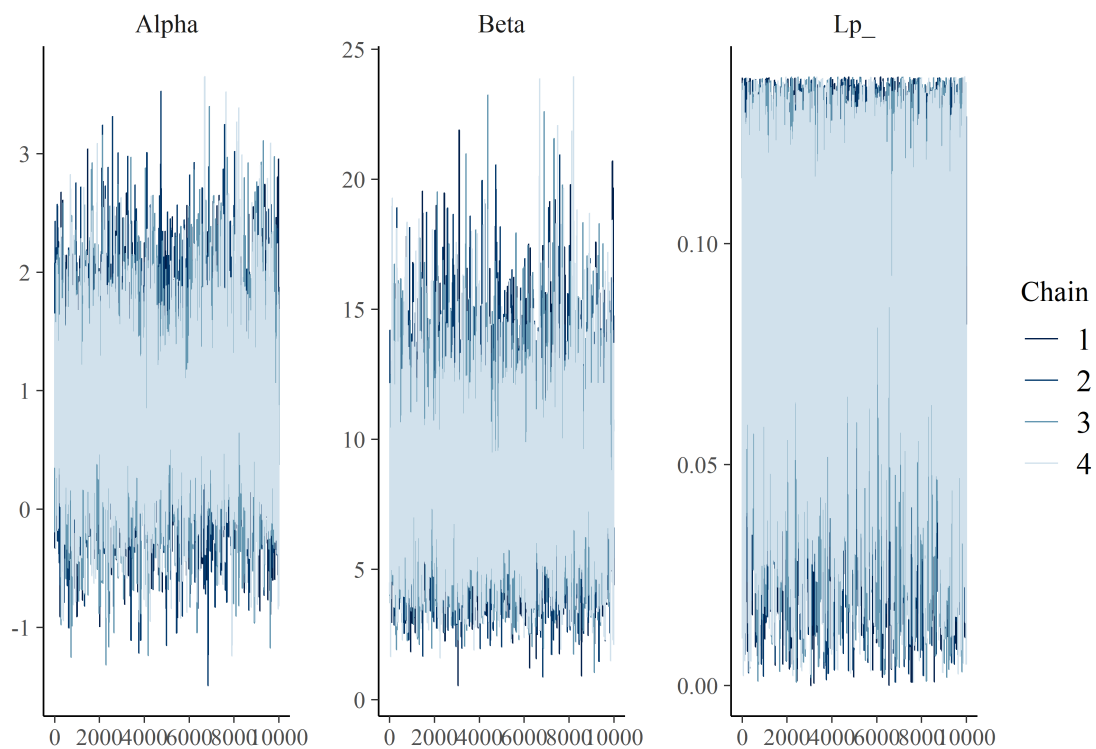
cs = c("Alpha","Beta","Lp_")
sims = mcmc_array(ns,nchains = nc, params = cs)
for(j in 1:nc){
  for(i in 1:ns){
    a.s = rnorm(n = 1,mean = a.c,sd = a.sd)
    b.s = rnorm(n = 1,mean = b.c,sd = b.sd)
```

```

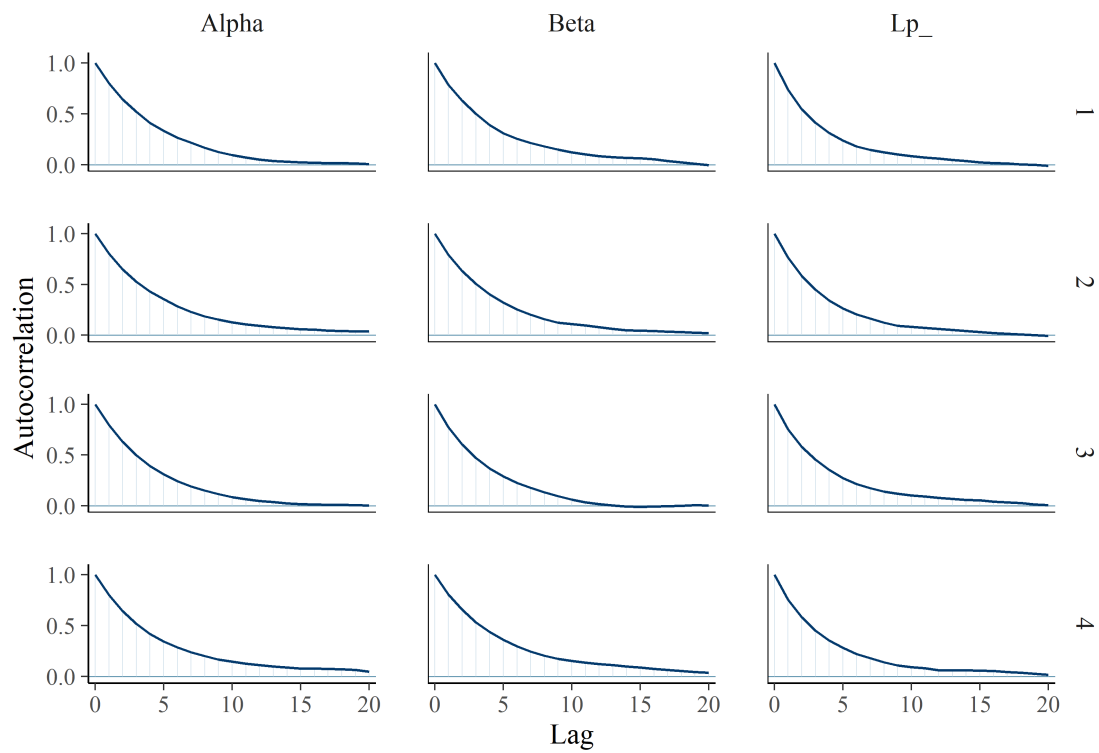
up = po(al = a.s,be = b.s)*dnorm(a.s,mean = 0.8,sd = 1)*dnorm(b.s,mean = 7.7,sd = 4.9)
do = po(al = a.c,be = b.c)*dnorm(a.c,mean = 0.8,sd = 1)*dnorm(b.c,mean = 7.7,sd = 4.9)
r = up/do
a = runif(1)
if(r > a){
  a.c = a.s
  b.c = b.s
}
post = po(al = a.c,be = b.c)
sims[i,j,] = c(a.c,b.c,post)
}
}

```

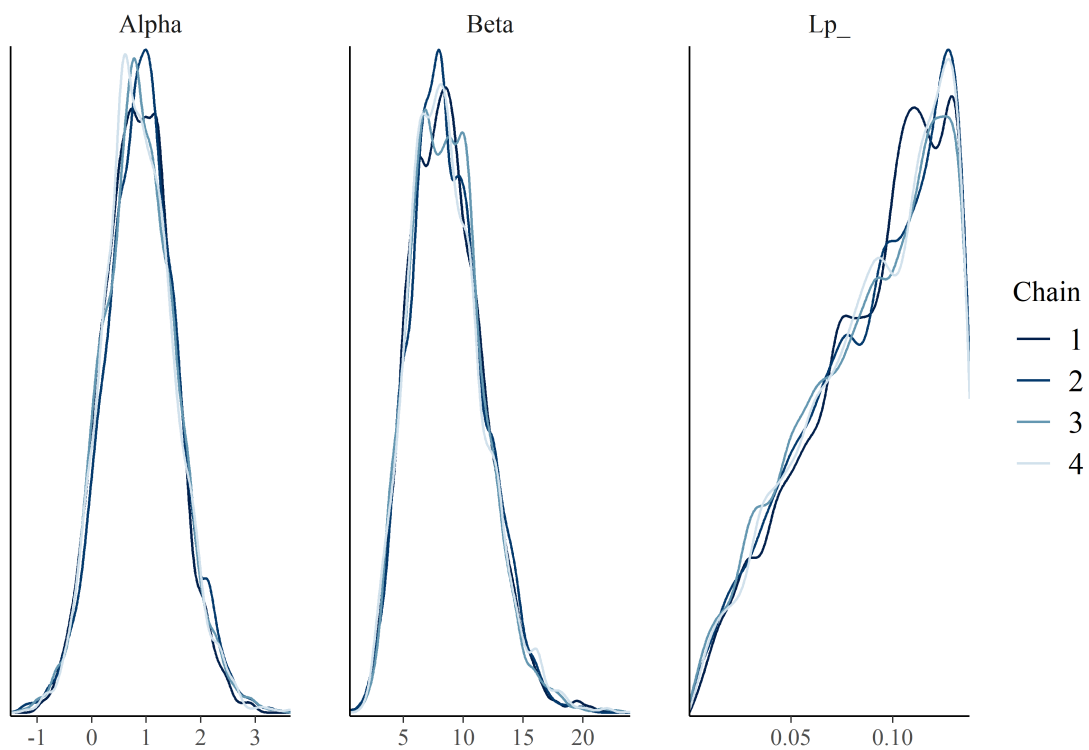
```
mcmc_trace(sims, cs)
```



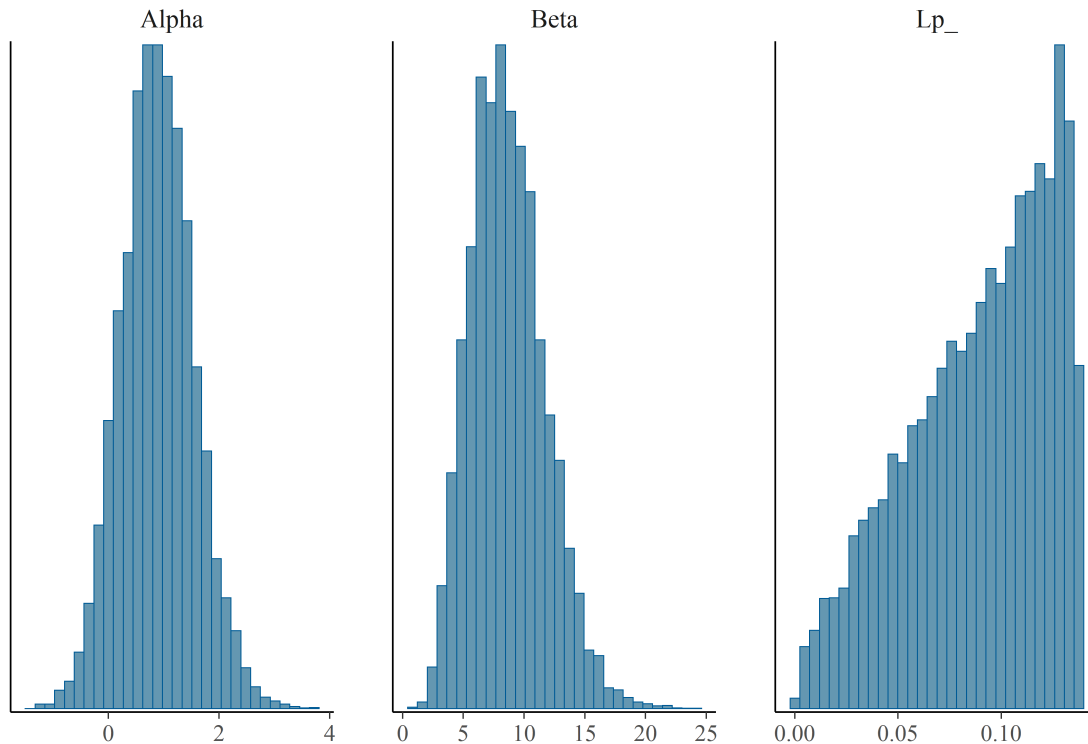
```
mcmc_acf(sims, cs)
```



```
mcmc_dens_overlay(sims, cs)
```



```
mcmc_hist(sims, cs)
```



BDA Problem 11.3

```
mo1 = stan_model("machine.stan")
```

```
monitor(sf)
```

```
## Inference for the input samples (4 chains: each with iter = 10000; warmup = 0):
```

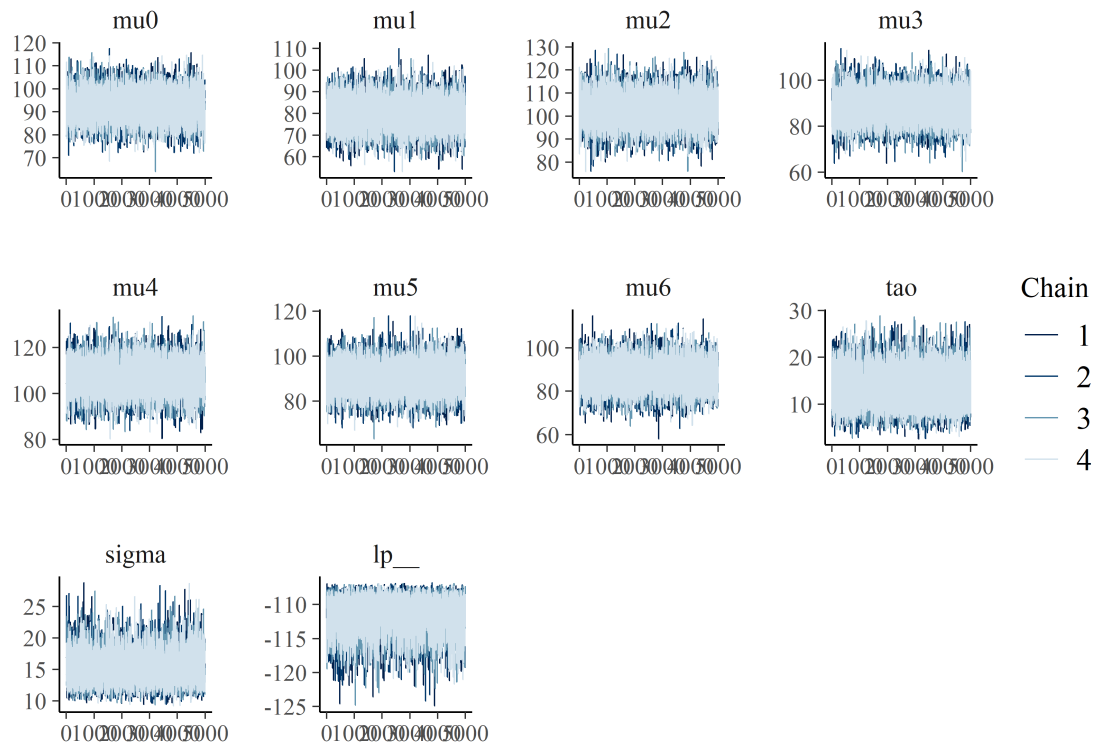
```
##
```

	Q5	Q50	Q95	Mean	SD	Rhat	Bulk_ESS	Tail_ESS
## mu0	83.4	93.0	102.5	93.0	5.8	1	17057	14873
## mu1	69.5	79.9	90.7	80.0	6.5	1	16549	10034
## mu2	92.5	103.2	113.3	103.1	6.3	1	21104	13797
## mu3	79.1	89.0	99.1	89.0	6.1	1	21704	13084
## mu4	96.4	107.3	117.9	107.2	6.5	1	17792	12815
## mu5	80.5	90.6	100.7	90.7	6.2	1	22011	13809
## mu6	77.5	87.6	97.6	87.6	6.1	1	19752	13250
## tao	7.6	13.2	19.9	13.4	3.8	1	12405	8199
## sigma	11.9	14.9	19.4	15.2	2.3	1	14983	10809
## lp__	-115.7	-111.0	-108.2	-111.4	2.3	1	7857	11462

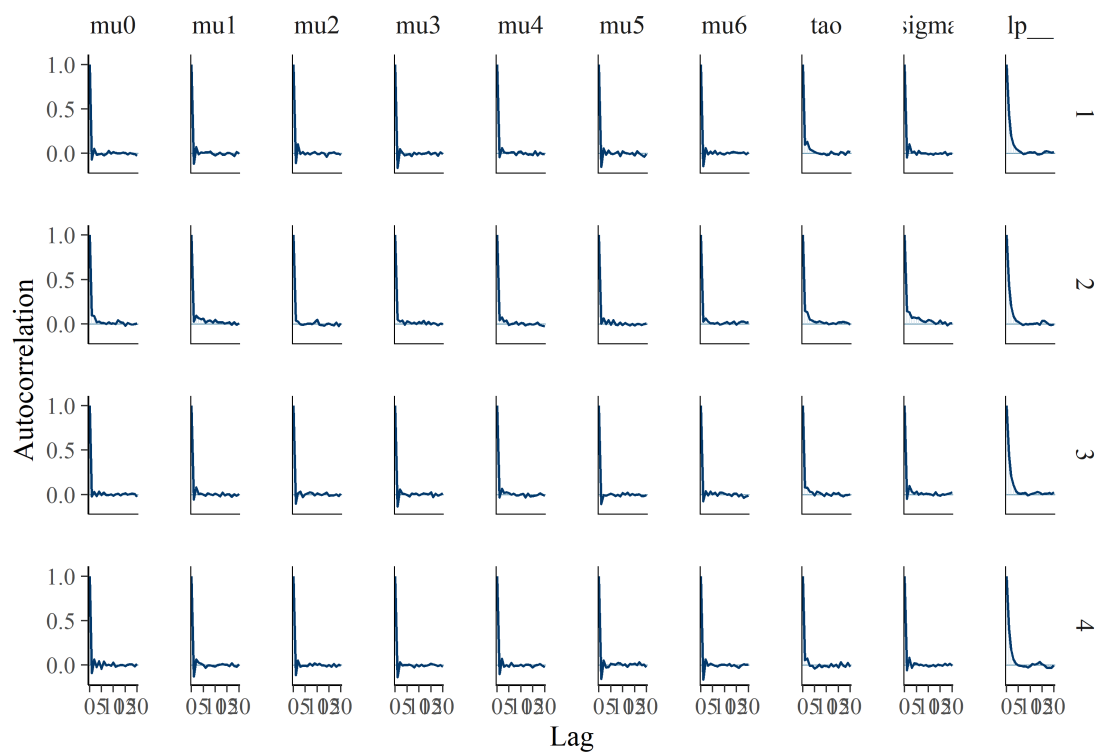
```
##
```

```
## For each parameter, Bulk_ESS and Tail_ESS are crude measures of
## effective sample size for bulk and tail quantities respectively (an ESS > 100
## per chain is considered good), and Rhat is the potential scale reduction
## factor on rank normalized split chains (at convergence, Rhat <= 1.05).
```

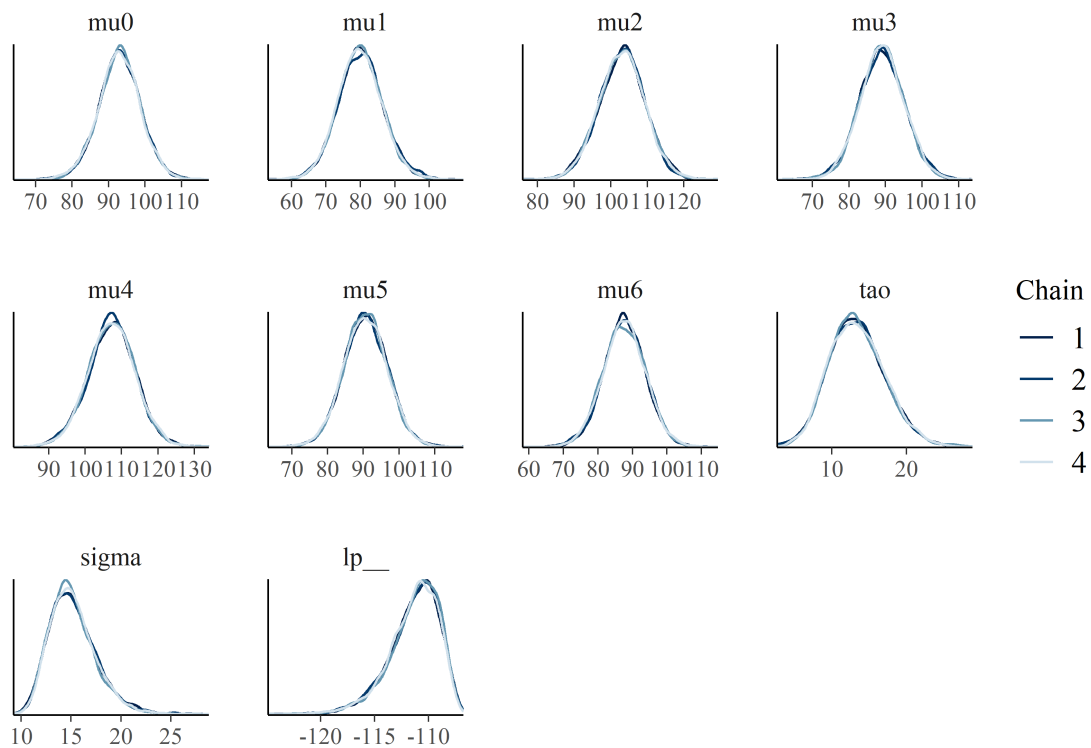
```
sims <- as.array(sf)
params <- c(paste0("mu", 0:6), "tao", "sigma", "lp__")
mcmc_trace(sims, params)
```



```
mcmc_acf(sims, params)
```



```
mcmc_dens_overlay(sims, params)
```



```
mcmc_hist(sims, params)
```

