

A combinator library for MCMC kernels and proposals

Praveen Narayanan, May 8 2014

Motivation

MCMC samplers are
complicated

$$\mathcal{A}(x^{(i)}, x^\star) = \min \left\{ 1, \frac{p(x^\star)q(x^{(i)} | x^\star)}{p(x^{(i)})q(x^\star | x^{(i)})} \right\}$$

$$\mathcal{A}_{n \rightarrow m} = \min \left\{ 1, \frac{p(m, x_m^*)}{p(n, x_n)} \times \frac{q(n | m)}{q(m | n)} \times \frac{q_{m \rightarrow n}(u_{m,n} | m, x_m^*)}{q_{n \rightarrow m}(u_{n,m} | n, x_n)} \times \mathcal{J}_{f_{n \rightarrow m}} \right\}$$

$$\mathcal{A}_{split} = \min \left\{ 1, \frac{p(k+1, \mu_{k+1})}{p(k, \mu_k)} \times \frac{\frac{1}{k+1}}{\frac{1}{k}} \times \frac{1}{p(u_{n,m})} \times \mathcal{J}_{split} \right\}$$

$$A_{merge} = \min \left\{ 1, \frac{p(k-1, \mu_{k-1})}{p(k, \mu_k)} \times \frac{\frac{1}{k-1}}{\frac{1}{k}} \times \mathcal{J}_{merge} \right\}$$

We would like
predefined samplers

MCMC depends on
“good” proposals

A I-D target distribution

$$p(x) \propto 0.3 \exp(-0.2x^2) + 0.7 \exp(-0.2(x - 10)^2)$$

**A gaussian might
work as a proposal**

**What is the optimal
mean? Standard
deviation?**

We would like to easily
experiment with **different**
proposals on a sampler

A second look

$$p(x) \propto 0.3 \exp(-0.2x^2) + 0.7 \exp(-0.2(x - 10)^2)$$

A mixture of
gaussians might
work even better

We would like to
easily **mix** various
proposals

Quick info

Proposal and target combinators

Kernel combinators

Making the random walk

Future work

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The library is
written in Haskell

**It extends Indiana's
probabilistic
programming system**

**Consists of three
modules:
Distributions,
Actions, Kernels**

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Targets vs proposals

Targets have
a density

type Density a = a → Probability

data Target a = T (Density a)

Proposals have a
density and can be
sampled from

```
type Sample a = Rand → IO a
```

```
data Proposal a =  
  P (Density a) (Sample a)
```

Proposals

**What proposals are
already defined?**

**uniform, normal,
categorical**

uniform [0,2,4] [3,5,7]

normal [0,1,4,7] (diag [2,2,2,2])

categorical

["pizza", "sushi", "curry"]

[0.7, 0.9, 0.3]

**How do we create
our own proposals?**

Mix existing
proposals

```
proposalMix ::  
[Proposal a] → [Double] → Proposal a
```

```
gMix = let g1 = normal [0,0] (diag [1,1])  
          g2 = normal [5,5] (diag [2,2])  
        in proposalMix [g1, g2] [0.3, 0.7]
```

**Increase dimensions
of existing proposals**

Use `updateNth`,
`updateBlock`

updateNth :: Int
→ ([a] → Proposal [a])
→ ([a] → Proposal [a])

**condProposal =
updateNth
4**

($\lambda y \rightarrow \text{normal } y \text{ } [[\mathbf{1}]]$)

Update component a_4
for a point $(a_1, a_2, a_3, a_4, \dots, a_n)$

updateBlock :: Int

→ Int

→ ([a] → Proposal [a])

→ ([a] → Proposal [a])

**condProposal =
updateBlock**

1

2

($\lambda y \rightarrow \text{normal } y \text{ (diag [1,1])}$)

Update components a_1, a_2
for a point (a_1, a_2, \dots, a_n)

Construct **our own**
using **makeProposal**

```
customProposal =  
let density = ...  
    sample = ...  
in makeProposal density sample
```

Targets

**How do we create
targets?**

Use `makeTarget`

```
t1 = makeTarget $  
    λ[x] → 0.3 * exp (-0.2*x*x)
```

Mix existing targets,
using **targetMix**

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**What is a
kernel?**

A **function** that,
when given a target
and conditional
proposal, performs
MCMC **steps**

type Step x = Rand → x → IO x

type Kernel x a =
Target a →
(a → Proposal a) →
Step x

**What kernels are
already defined?**

metropolisHastings,
simulatedAnnealing

How do we define
our own steps?

Mix some
steps

mixSteps :: [Step x]
→ [Double]
→ Step x

```
mhMix = let mh = metropolisHastings
          mh1 = mh target proposal1
          mh2 = mh target proposal2
        in mixSteps [mh1, mh2] [0.7, 0.3]
```


Cycle

through a
kernel

cycleKernel :: Kernel x a
→ Target a
→ [a → Proposal a]
→ Step x

```
mhCycle =  
cycleKernel  
metropolisHastings  
target  
[p1, p2, p3]
```

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walk :: Step x

→ **X**

→ **Int**

→ **Rand**

→ **Action x a**

→ **IO a**

Start state

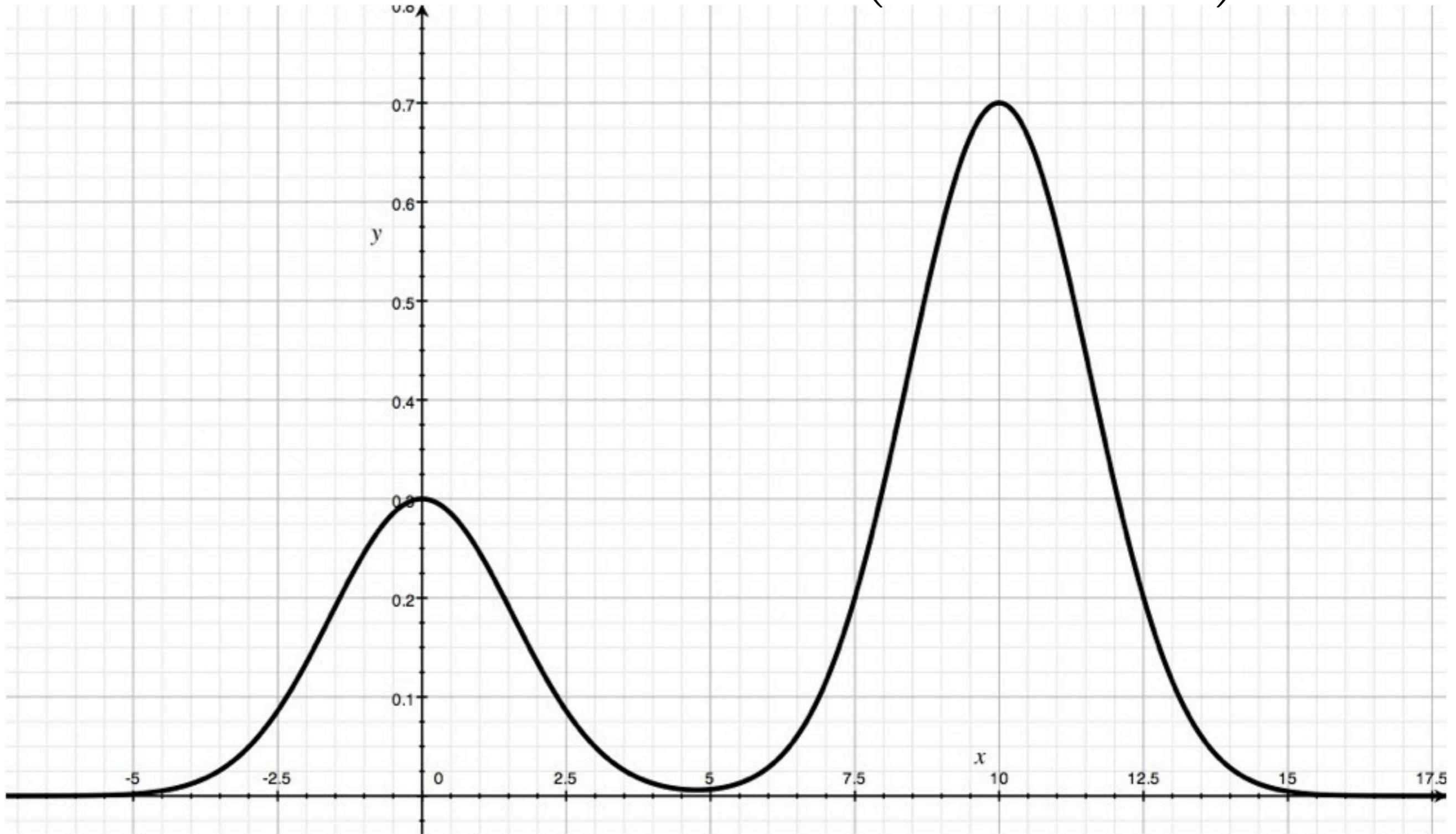
Number of steps

Action to take at each step

A third look

$$p(x) \propto 0.3 \exp(-0.2x^2) + 0.7 \exp(-0.2(x - 10)^2)$$

$$0.3 \exp(-0.2x^2) + 0.7 \exp(-0.2(x-10)^2)$$



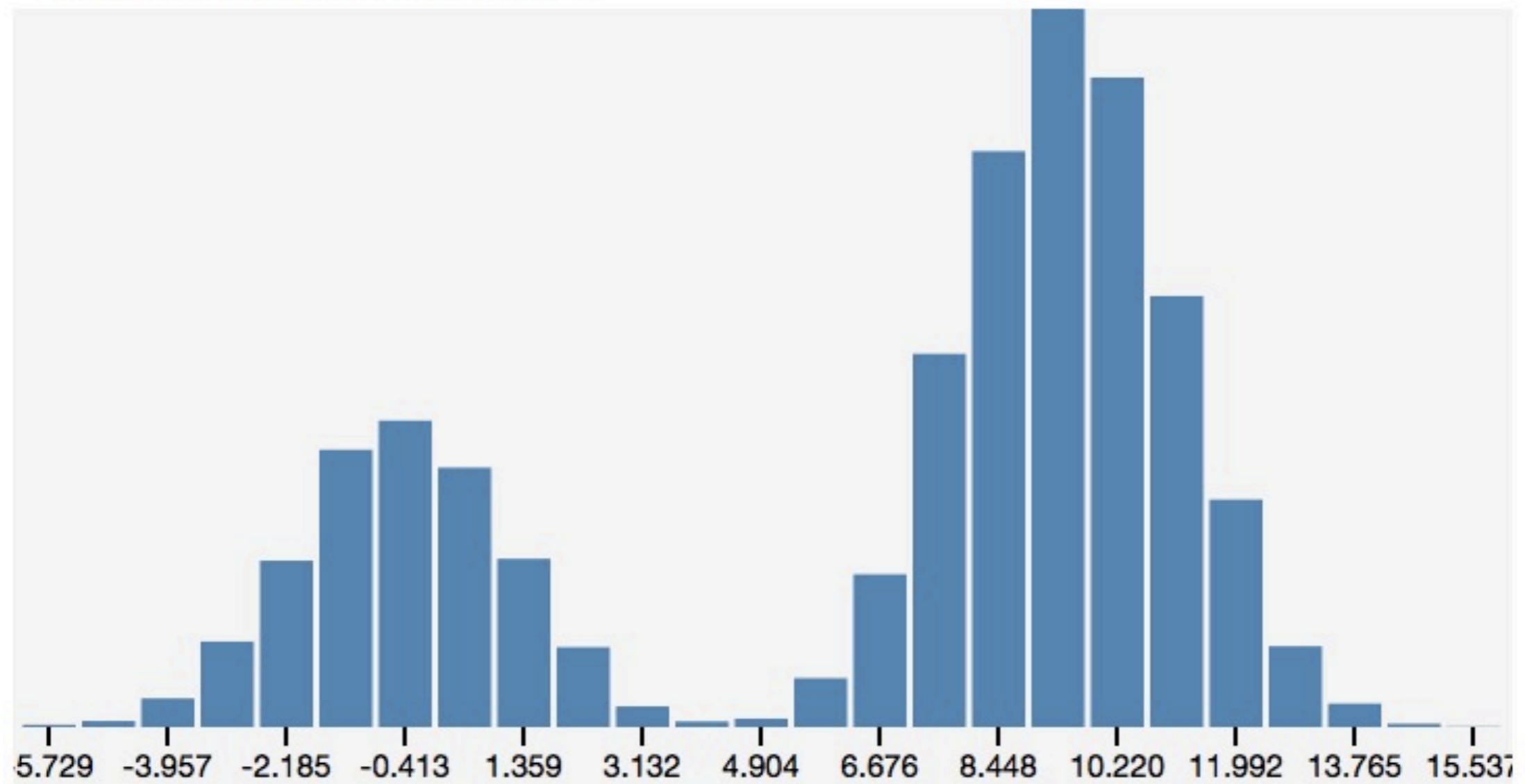
```
mhTest :: IO ()
mhTest = do
  g ← MWC.createSystemRandom
  let (t,p) = (... , ...)
      mh = metropolisHastings t p
      act = every 100 (batchViz vizMH 50)
  walk mh [0] (106) g act
```


Number of samples plotted: 10000

x

Min: -5.7291215850716135

Max: 16.422778145776434



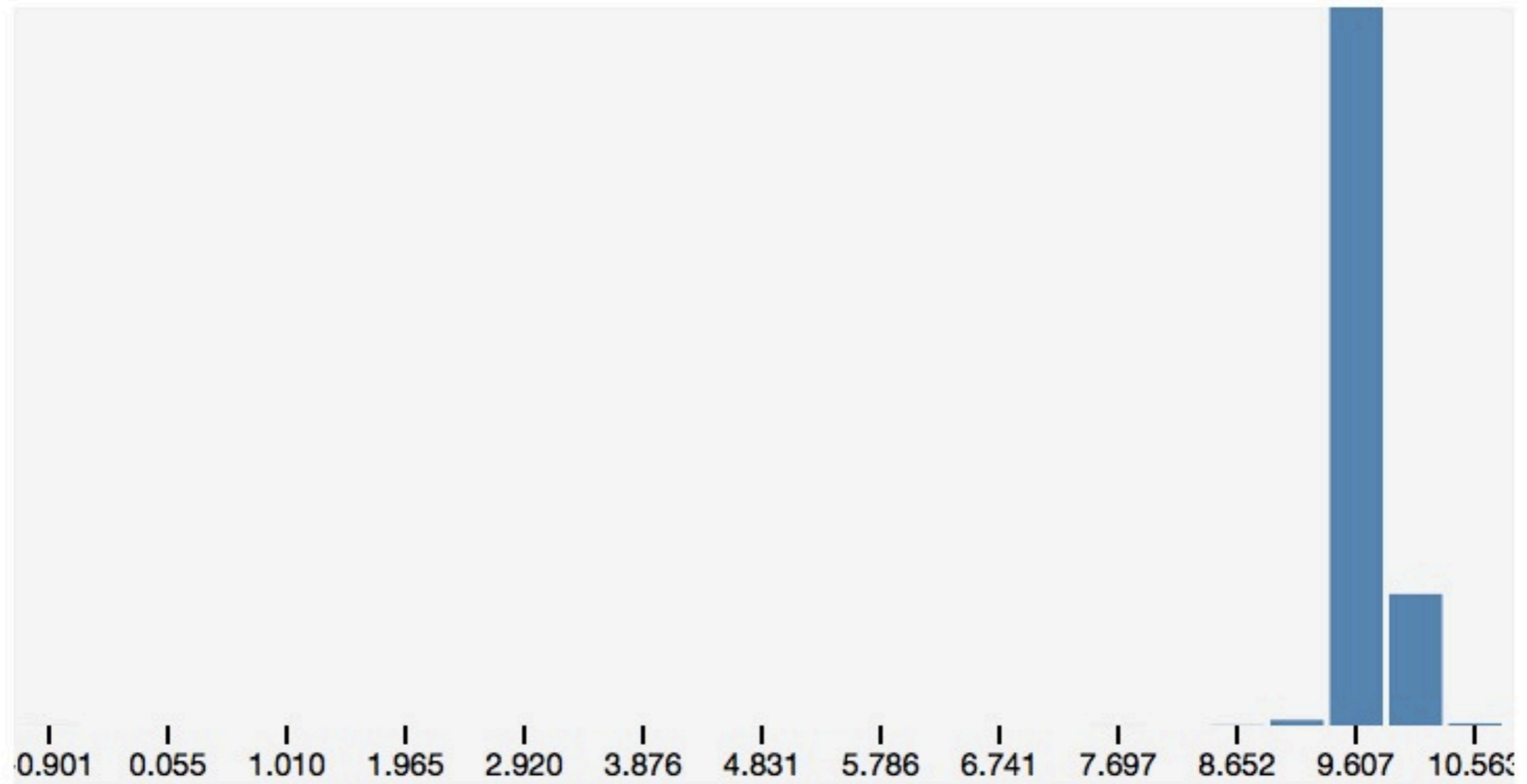
```
saTest :: IO ()
saTest = do
  g ← MWC.createSystemRandom
  let (t,p) = (... , ...)
      sa = simulatedAnnealing t p
      coolSch temp = temp * (1 - 1e-3)
      x0 = ([0], 1, coolSch)
      act = every 100 (batchViz vizSA 50)
  walk sa x0 (10^6) g act
```

Number of samples plotted: 9889

X

Min: -0.9005223518480108

Max: 11.04013597514839



```
blockMH :: Step [Double]
```

```
blockMH = let target = normal [0,1,4,7] (diag [2,2,2,2])  
           mh4D = metropolisHastings target  
           mh1 = mh4D $ updateBlock 1 2  
                   (λy → normal y (diag [1,1]))  
           mh2 = mh4D $ updateBlock 3 3 (λy → normal y [[1]])  
           mh3 = mh4D $ updateNth 4 (λy → normal y [[5]])  
       in mixSteps [mh1, mh2, mh3] [0.5, 0.4, 0.7]
```

```
blockTest :: IO ()
blockTest = do
  g ← MWC.createSystemRandom
  let a = batchPrint printMH 50
  walk blockMH [0,0,0,0] (106) g a
```

[0.907,-1.067,2.128,5.558],
[0.907,-1.067,2.128,6.584],
[1.336,-1.198,2.128,6.584],
[1.336,-1.198,2.128,6.584],
[-0.857,-1.729,2.128,6.584],
[-0.857,-1.729,2.660,6.584],
[-0.857,-1.729,2.660,6.584],
[-0.857,-1.729,2.660,8.325],
[-0.857,-1.729,2.660,8.325],
[-0.857,-1.729,2.660,8.325],
[0.598,-2.241,2.660,8.325],
[0.598,-2.241,2.660,8.325],
[0.598,-2.241,2.660,8.325],
[0.598,-2.241,2.660,7.699],
[0.598,-2.241,2.660,7.699], ...

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**Kernel combinators
for changing the
accept ratio**

More sampling methods

Approximate MH,
Stochastic EM,
Reversible jump
MCMC

Thank you