# Problemset 1

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### 1.1 Random Variable Generation

### a) Implementation of Lehmer RNG

We implement the Lehmer RNG:

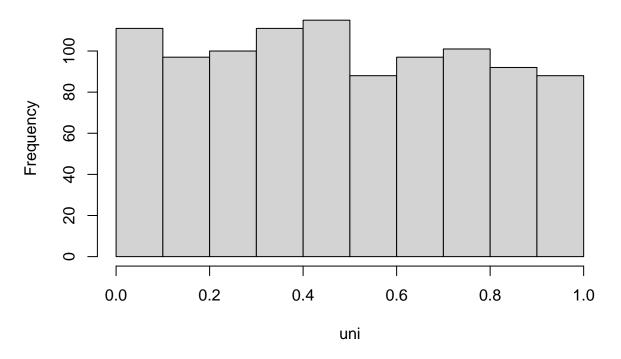
$$x_n = (ax_{n-1}) \bmod m$$

The seed  $x_0$  has to be chosen beforehand,  $m=2^31-1,\,a=7^5$ 

```
LRNG <- function(n, seed = 42) {
    m <- 2^31 -1 # .Machine$integer.max
    a <- 7^5
    samp <- numeric(n)
    samp[[1]] <- (a * seed) %% m
    for (i in seq_len(n - 1)) {
        samp[[i + 1]] <- (a * samp[[i]]) %% m
    }
    return(samp / m)
}</pre>
```

We simulate n = 1000 samples and plot as a histogram:

# Histogram of uni



### b) Implement inverse transformation method

```
ITM <- function(inverse, n, seed = 42) {
  uni <- LRNG(n = n, seed = seed)
  inverse(uni)
}</pre>
```

We obtain the inverse:

$$F(x) = 1 - e^{-2x} (1)$$

$$\Longrightarrow F^{-1}(q) = -0.5 * log(1-q) \tag{2}$$

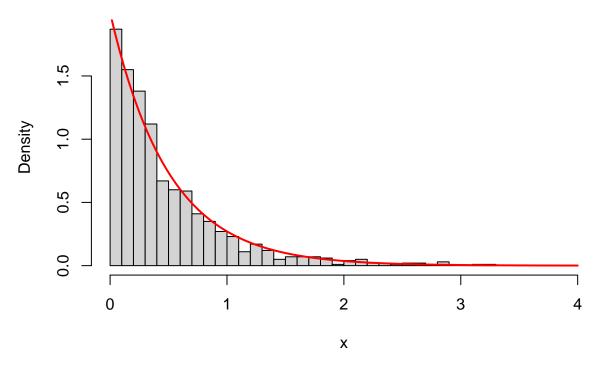
Subsequently we call ITM on the inverse. Additionally we implement the true density  $f_X(x) = 2e^{-2x}$ .

```
Fq <- function(q) {
    -.5 * log(1 - q)
}

fx_samp <- ITM(Fq, 1000)
fx_true <- function(x) {2*exp(-2*x)}</pre>
```

We plot the true density vs. the simulated samples:

## sample vs true density



### c) Goodness of fit testing

 ${\bf Kolmogorov\text{-}Smirnov\text{-}Test}$ 

```
# ?ks.test
ks.test(fx_samp, "pexp", rate = 2)
##
##
   Asymptotic one-sample Kolmogorov-Smirnov test
## data: fx_samp
## D = 0.044486, p-value = 0.0382
## alternative hypothesis: two-sided
\mathcal{X}^2-Test
chisq <- function(samp, bins = 10, pdf, ...) {</pre>
  tab <- table(cut(c(0, samp), breaks = bins)[2:(length(fx_samp)+1)])</pre>
  step \leftarrow max(samp) / 10
  p <- vapply(seq_len(bins),</pre>
               function(x) pdf(x*step, ...) - pdf((x - 1)*step, ...),
               numeric(1))
  chisq.test(tab, p = p, rescale.p = TRUE)
x2 <- chisq(fx_samp, pdf = pexp, rate = 2)</pre>
x2$p.value
## [1] 0.2243432
```

#### x2\$observed ## (-0.00322, 0.322](0.322, 0.645](0.645, 0.967](0.967, 1.29]## ## 512 133 ## (1.29, 1.61](1.61, 1.93](1.93, 2.26](2.26, 2.58]## 26 20 ## (2.58, 2.9](2.9, 3.23]## x2\$expected ## (-0.00322,0.322] (0.322, 0.645](0.645, 0.967](0.967, 1.29]475.881025 249.777651 68.811964 ## 131.101833 ## (1.29, 1.61](1.61, 1.93](1.93, 2.26](2.26, 2.58]## 36.117621 18.957206 9.950147 5.222575 (2.58, 2.9]## (2.9, 3.23]2.741195 1.438782

Since the the  $\mathcal{X}^2$ -Test uses binning, it may be able to smooth out little deviations from the expectation. We would expect a higher p-value for the Kolmogorov-Smirnov-Test, if we would have simulated more samples.

```
ks.test(ITM(Fq, 100000), "pexp", rate = 2)

##
## Asymptotic one-sample Kolmogorov-Smirnov test
##
## data: ITM(Fq, 1e+05)
## D = 0.0033039, p-value = 0.225
## alternative hypothesis: two-sided
```

# 1.2 Rejection Sampling, Importance Sampling

### a) Rejection Sampler Implementation

The maximum of  $exp(-x^2)$  is located at  $x_0 = 0$ , with a value of exp(0) = 1, therefore we need to scale our umbrella-distribution to be  $\geq 1$ .

Since the density of the Uniform-distribution an [-a,a] is  $\frac{1}{2a}$ , we search for  $\alpha$  such that  $\alpha \cdot u \geq 1 \implies \alpha \geq 2a$ , where  $u \sim U(-a,a)$ .

```
rejection_sampler <- function(n_samples, a, verbose = FALSE, print.steps = FALSE) {
   checkmate::assertIntegerish(n_samples, lower = 1, len = 1, any.missing = FALSE)
   checkmate::assertNumeric(a, lower = .Machine$double.eps, len = 1, any.missing = FALSE)
   checkmate::assertFlag(verbose)
   checkmate::assertFlag(print.steps)

###
   alpha <- 2*a + .Machine$double.eps
   f <- function(x) exp(-x^2) * (x <= a) * (x >= -a)
   g <- function(x) (1 / 2*a) * (x <= a) * (x >= -a)
   G_inv <- function(q) 2*a*q - a

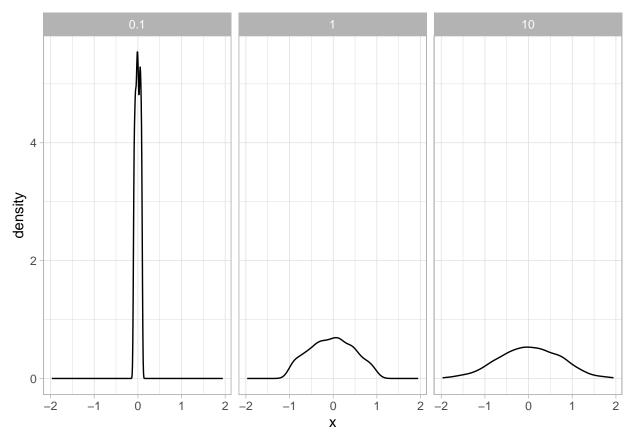
###
   samp <- numeric(n_samples)
   start <- 1
   step <- 1
   while(n_samples > 0) {
```

```
Y <- G_inv(runif(n_samples))
   U <- runif(n_samples)</pre>
   test <- U \leftarrow f(Y) / (alpha * g(Y))
   approved <- sum(test)</pre>
   if(verbose) {
      cat(sprintf("step %d: approved %d of %d remaining samples...\n",
                 step, approved, n_samples))
   step <- step + 1
   if (approved > 0) {
     samp[start:(start + approved - 1)] <- Y[test]</pre>
     start <- start + approved
     n_samples <- n_samples - approved
  }
  if(print.steps) {
   cat(sprintf("#########"\n number of necessary steps: %d \n#########\n",
                 step))
 }
  return(structure(samp, steps = step)
}
head(rejection_sampler(n = 1000, a = 1, verbose = TRUE))
## step 1: approved 743 of 1000 remaining samples...
## step 2: approved 189 of 257 remaining samples...
## step 3: approved 51 of 68 remaining samples...
## step 4: approved 15 of 17 remaining samples...
## step 5: approved 2 of 2 remaining samples...
attr(rejection_sampler(n = 1000, a = 1), "steps")
## [1] 6
```

### b) Sampling for different values of a

```
We sample n = 1000 with a \in \{0.1, 1, 10\}
```

```
tibble(a = c(.1, 1, 10)) %>%
mutate(x = lapply(a, function(a) rejection_sampler(1000, a))) %>%
unnest_longer(x) %>%
ggplot(aes(x)) +
geom_density() +
facet_wrap(~a, ncol = 3) +
theme_light()
```



With bigger a, we see that we have a bigger variance in the cdf.

### c) Importance Sampling

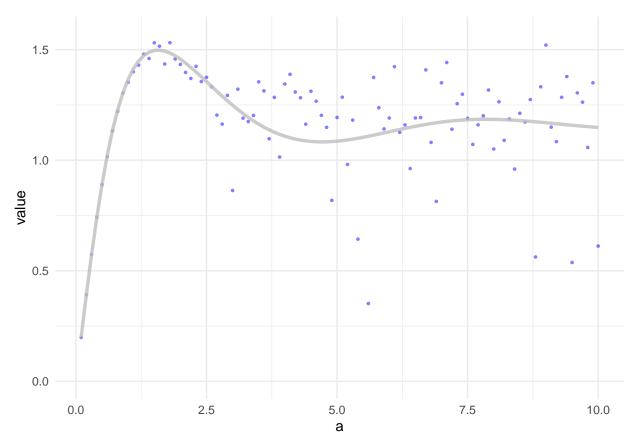
We want to estimate  $I_a = \int_{-a}^a \frac{\cos(x)}{1+x^2} dx$ . Upon multiplying with  $\frac{g(x)}{g(x)} = \frac{\exp(-x^2)}{\exp(-x^2)}$  we obtain  $I_a = \mathbb{E}_G(h(X)\frac{f(X)}{g(X)}) = \int_{-a}^a \frac{1}{1+x^2} \frac{\cos(x)}{\exp(-x^2)} \exp(-x^2) dx$ .

```
integrand <- function(x) cos(x)/(1 + x^2)

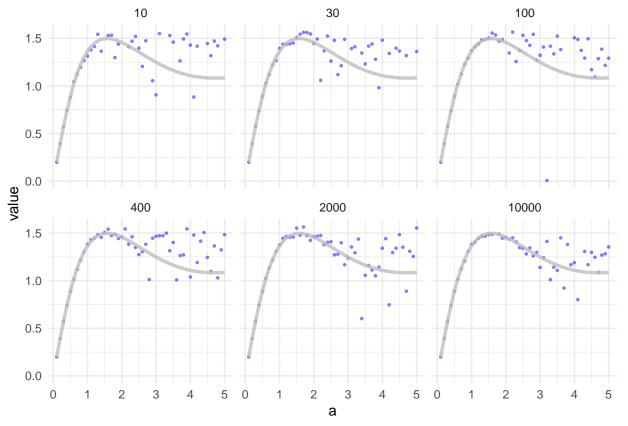
importance_sampler <- function(n = 1000, a = 1) {
    g <- function(x) exp(-x^2)
    ### calculate normalization constant
    nu <- mean(g(runif(1000, -a, a))) * 2 * a
    Y <- rejection_sampler(n_samples = n, a = a)
    mean(integrand(Y) / g(Y) * nu)
}

### sampler2: use density of continious uniform function on [-a,a]
importance_sampler2 <- function(n = 1000, a = 1) {
    U <- runif(1000, -a, a)
    ### calculate normalization constant
    nu <- 2*a
    mean(integrand(U) * 2 * a)
}</pre>
```

Plot the samples (blue) for different values of a. The grey line is the numerical solution to the integral.



Show behaviour for different values of n:



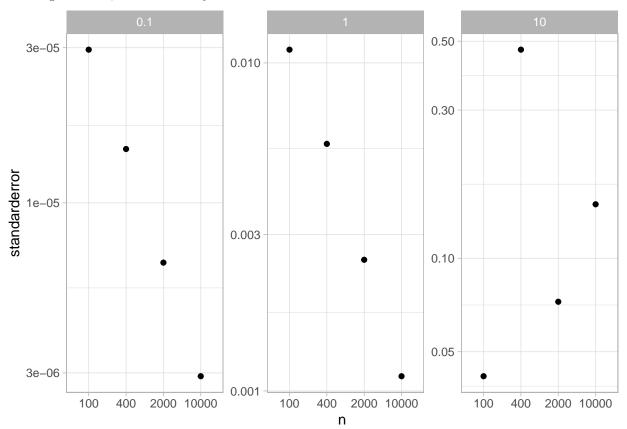
## d) Computing standard error

We compute the standard error as  $var[I_a] = var[\mathbb{E}_G(h(X)\frac{f(X)}{g(X)})] = \frac{var[h(X)\frac{f(X)}{g(X)}]}{n}$ 

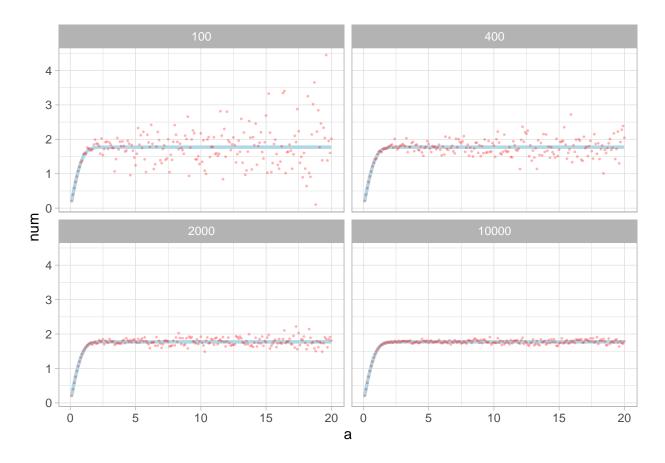
```
se_est <- function(n, a) {
   g <- function(x) exp(-x^2)
   integrand <- function(x) cos(x)/(1 + x^2)

### calculate normalization constant
   nu <- mean(g(runif(1000, -a, a))) * 2 * a
   Y <- rejection_sampler(n_samples = n, a = a)
   sigma <- sqrt(var(integrand(Y) / g(Y) * nu) / n)
   return(sigma)
}</pre>
```

Plotting se over a, for different n yields:



Plotting the samples for different n over a vs. the expected value:



## e) Can this method be used for $a \to \infty$ ?

By using very large values for a, the amount of rejected samples is increasing sharply.

The amount of steps needed to compute the rejection sample for incresing values of a are shown below. The number of steps can be interpreted as complexity of the operation. It can be visually approximated that  $steps = \mathcal{O}(a^3), a \in [0, 20]$ 

