

# 1 Mathematical preparations

## 1.1 f(x)

pdf

$$\begin{aligned} f(x) &= e^{-x} \\ x &\in (0, \infty) \end{aligned}$$

normalization:

$$\begin{aligned} \int_0^\infty f(x) dx &= \int_0^\infty e^{-x} dx \\ &= -e^{-x} \Big|_0^\infty \\ &= 1 \end{aligned}$$

cdf

$$\begin{aligned} F(x) &= \int_0^x f(t) dt \\ &= \int_0^x e^{-t} dt \\ &= -e^{-t} \Big|_0^x \\ &= -e^{-x} + 1 \end{aligned}$$

invert:

$$\begin{aligned} y &= -e^{-x} + 1 \\ \ln(1 - y) &= -x \\ F^{-1}(y) &= -\ln(1 - y) = x \end{aligned}$$

## 1.2 g(x)

pdf

$$\begin{aligned} g(x) &= x \cdot e^{-\frac{x^2}{\alpha^2}}, \quad \alpha = 25^\circ \\ x &\in [0^\circ, 180^\circ] \end{aligned}$$

normalization:

$$\int_{0^\circ}^{180^\circ} g(x) dx = \int_{0^\circ}^{180^\circ} x \cdot e^{-\frac{x^2}{\alpha^2}} dx$$

substitute  $\frac{x^2}{\alpha^2} =: b$ ,  $x \cdot dx = \frac{\alpha^2}{2} db$

$$\int_{0^\circ}^{180^\circ} g(x) dx = \frac{\alpha^2}{2} \int_{0^\circ}^{(180^\circ/\alpha)^2} e^{-b} db$$

$$\begin{aligned}
&= -\frac{\alpha^2}{2} [e^{-b}]_0^{(180^\circ/\alpha)^2} \\
&\approx \frac{\alpha^2}{2}
\end{aligned}$$

redefine  $g(x)$ :

$$g(x) = \frac{2}{\alpha^2} x \cdot e^{-\frac{x^2}{625\sigma}}$$

cdf

$$\begin{aligned}
G(x) &= \int_0^x g(t) dt \\
&= \frac{2}{\alpha^2} \int_0^x e^{-\frac{t^2}{625\sigma}} dt
\end{aligned}$$

substitute  $\frac{t^2}{\alpha^2} =: c$ ,  $dt = \frac{\alpha^2}{2} dc$

$$\begin{aligned}
G(x) &= \int_0^{x^2/\alpha^2} e^{-c} dc \\
&= -[e^{-c}]_0^{x^2/\alpha^2} \\
&= -e^{-x^2/\alpha^2} + 1
\end{aligned}$$

invert

$$\begin{aligned}
y &= -e^{-x^2/\alpha^2} + 1 \\
\ln(1-y) &= -\frac{x^2}{\alpha^2} \\
x &= \pm \alpha \sqrt{-\ln(1-y)} \\
G^{-1}(y) &= \pm \alpha \sqrt{-\ln(1-y)}
\end{aligned}$$

## 2 Computational Setup

For implementing the scattering of particles computed by a Monte Carlo method, 6 non-main methods were coded in addition to the main method.

The main-method is used only for initializing the random seed and starting the needed simulation.

For simulating, two different methods were implemented: `simulate_stat()` is used for getting the statistic properties of the scattering processes and therefore needs a number of particles that are computed. `simulate_path()` computes the path of 5 particles and saves it.

The process for simulating a particle path is the same in both methods: The path is computed by `getPath()`, if the particle becomes absorbed by `becomesAbsorbed()` and if it wasn't absorbed the scattering angle by `getAngle()`.

After computation the statistics are saved using `write()`.

When computing the path-length, the x-axis is spread by factor 10 to ensure greater accuracy. Therefore, the number of events for a certain path length has to be increased by factor 10 when being compared with  $f(x)$ .

When computing the scattering angle, via a random number the decision for scattering in positive or negative y-direction is made. Due to this, only half of the expected (by  $g(x)$ ) particles are measured with a certain scattering angle. Therefore, the statistics are not being compared to  $g(x)$  but to  $|\frac{g(x)}{2}|$ .

### 3 Observations

#### 3.1 Trajectories

The particle trajectories show that the particles become absorbed rather quickly. None of the displayed trajectories enters further into the medium than 6[mm].

#### 3.2 Path length

The path length distribution fits  $f(x)$  nicely. As expected, the variance between  $f(x)$  and the simulated distribution decreases with increasing number of simulated particles.

A bit strange is the behaviour of the path length for  $r = 0$ : Here, the measured number of particles stay way below (at about half) the expected value. This holds for all tested number of particles.

#### 3.3 Absorption ratio

The measured absorption ratio is about the exact value of 0.2. With increasing number of particles, again, the variance between expectation and measured values decreases.

#### 3.4 Scattering angle

The measured scattering angle is distributed almost exactly as given by  $g(x)$ . As in the above cases, the variance between measurement and expectation decreases with increasing number of sampled particles.

#### 3.5 Detector at $x=1$ and $x=5$

For a detector at  $x=1$  or  $x=5$ , a large maximum appears at  $y = 0$  and two side-maxima at  $y \approx \pm 0.25$ . This effect probably takes place because most events occur directly after inserting the particle at  $(0/0)$  with movement exactly along the x-axis. The side-maxima appear due to scattering processes and are therefore distributed as  $g(x)$ .

The scheme of maxima and side-maxima is clearer to be seen when the number

of simulated particles is increased.

It is no surprise that there are more events detected in a detector at  $x=1$  than at  $x=5$  (for  $N = 100$  even no event at all was registered at  $x=5$ ) since the second detector is further away from the point of insertion.