

# Diffusion

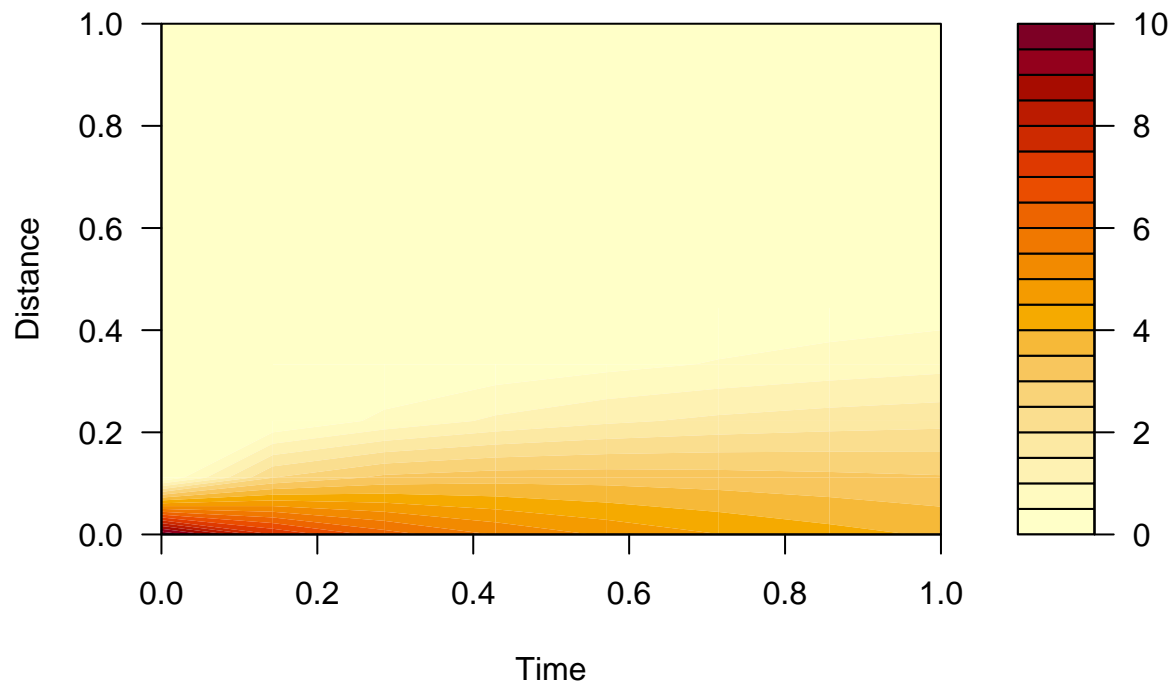
## R implementation

```
source(here("R","diffusion.R"))

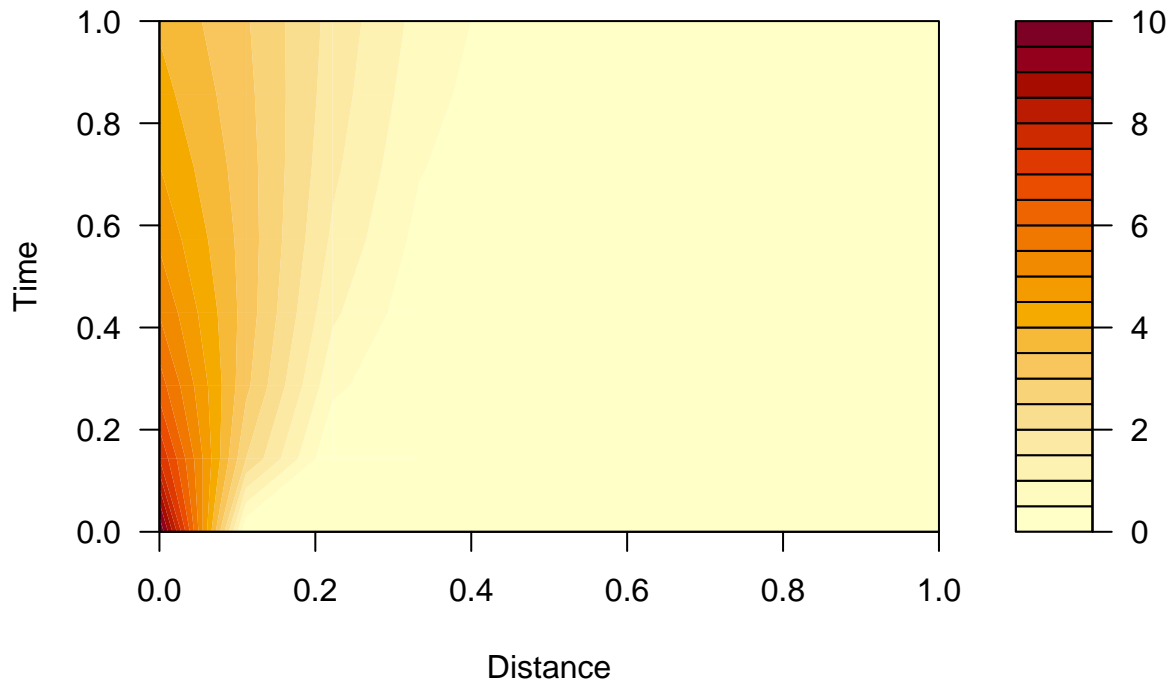
# run our diffusion model (iterative difference equation) with initial concentration of 10, for 8 times
# using diffusion parameters 0.5 s/m2, 10 m2
result = diff1(initialC=10, nx=10, dx=1, nt=8, dt=1, D=0.5, area=10)

# a list is returned with our 3 data frames for concentration (conc), qin and qout
# result

# used filled contour to plot results
# head(result$conc)
filled.contour(result$conc, xlab="Time", ylab="Distance")
```



```
# or if you prefer this orientation (Distance on x axis)
filled.contour(t(result$conc), ylab="Time", xlab="Distance")
```

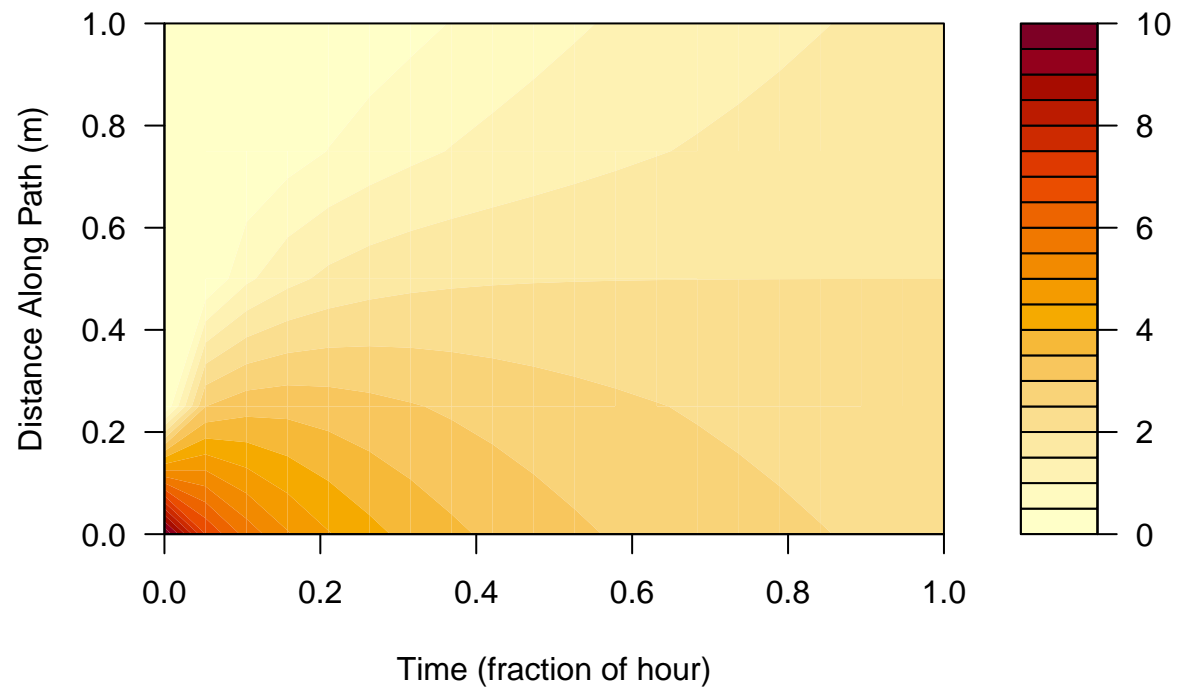


Change parameters (diffusivity  $D$ , and space and time steps ( $dx$ ,  $dt$ ))

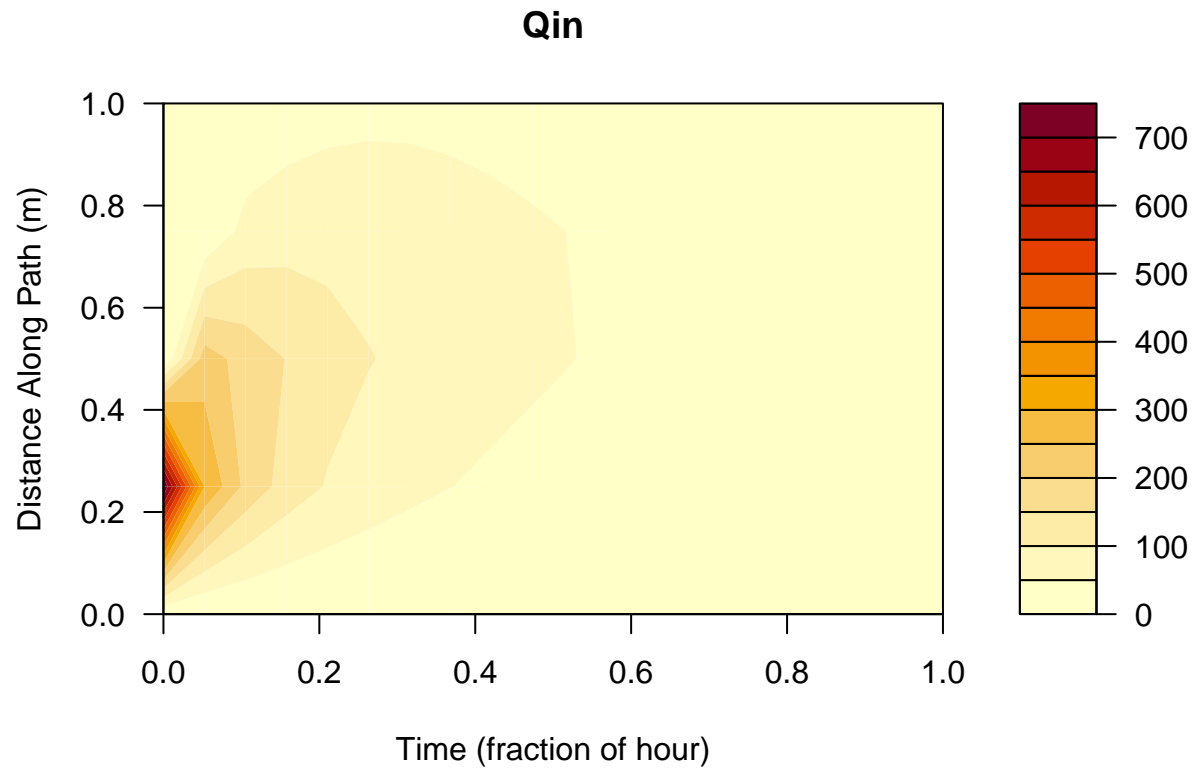
```
# changes diffusivity and other parameters particularly
# diffusivity, dx and dt
res=diff1(initialC=10,nx=5,dx=25,nt=20,dt=20,D=0.75,area=10)

filled.contour(res$conc, xlab="Time (fraction of hour)",ylab="Distance Along Path (m)", main="Pollutant
```

## Pollutant Diffusion



```
# we can also see how much material moved from place to place each time step  
filled.contour(res$qin, xlab="Time (fraction of hour)", ylab="Distance Along Path (m)", main = "Qin")
```



Overall, increasing the total simulation time ( $nt * dt$ ) lead to an increase in pollutant diffusion over time and space. Decreasing the number of discrete segments ( $m$ ) lead to greater distance of pollutants along the path.