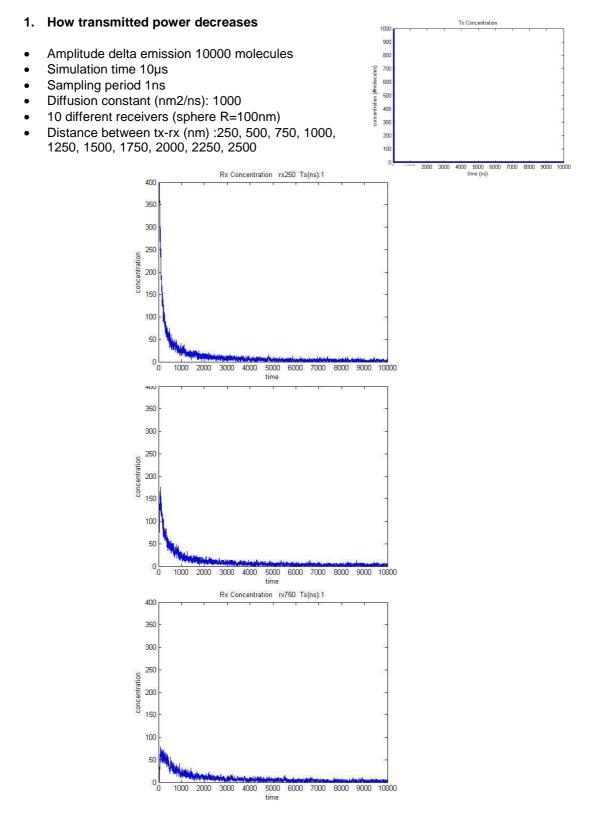
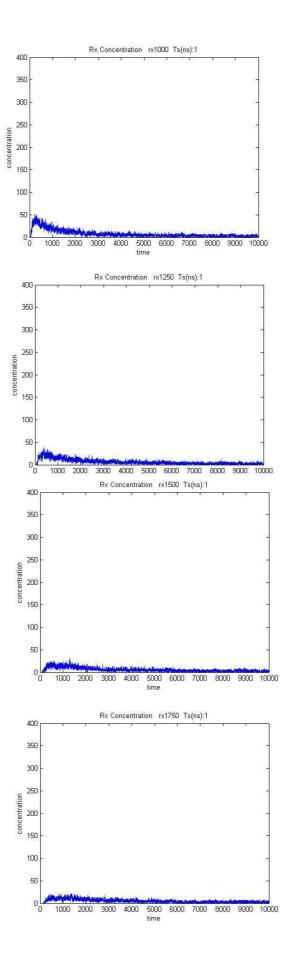
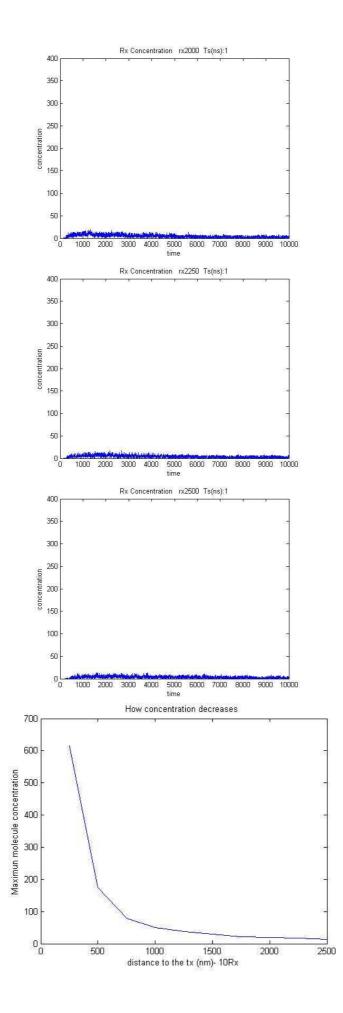
Simulations

These simulations are made with the first version of the diffusive molecular channel simulator, which implements an infinite space; we are simulating the channel without walls. The transmitter emits a specific number of molecules during a time and these molecules propagate due to Brownian motion until the receiver. This one measures the number of molecules within a specific volume. As there is no space limits we are not taking into account a possible initial concentration of molecules.

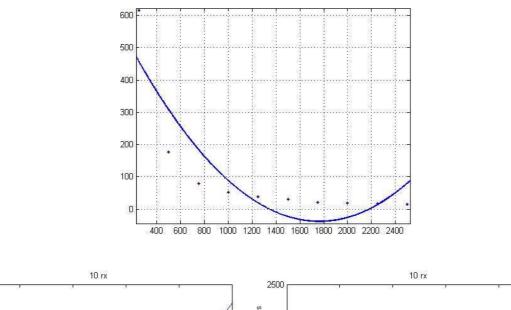


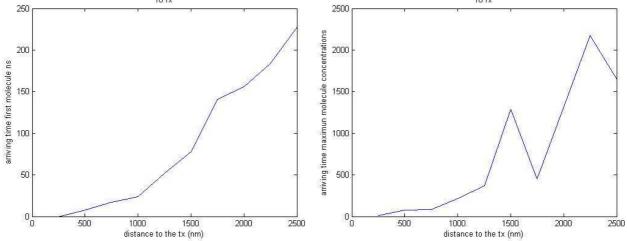


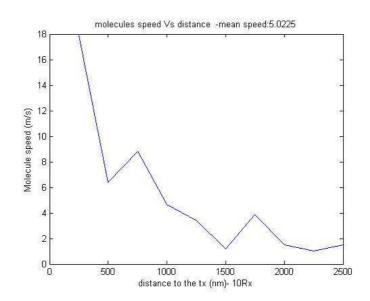


The concentration decrease fits with the following quadratic polynomial with a 95% confidence bounds: $f(x) = 0.000217 x^2 - 0.7646 x + 635.4$

- p1 = (5.499e-005, 0.000379)
- p2 = (-1.222, -0.3076)
- p3 = (361.8, 909)

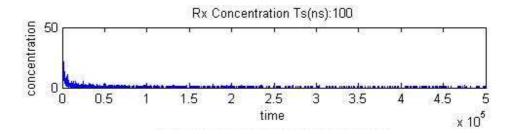


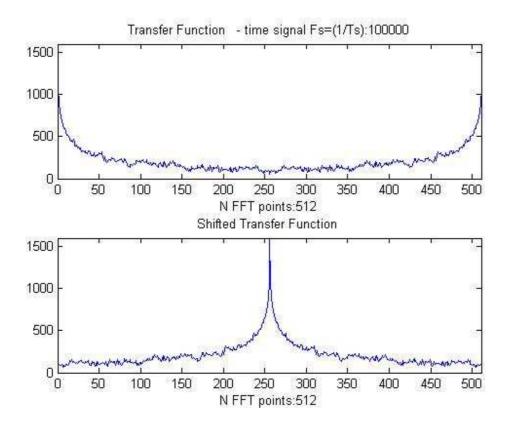




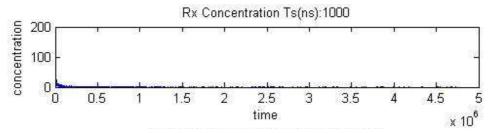
2. Channel Transfer Function

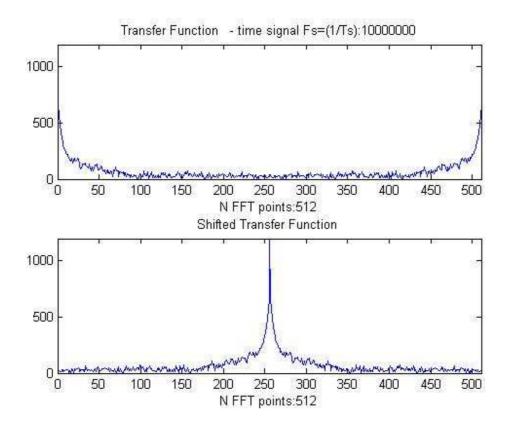
- Simulation time 500 µs
- Sampling Period 100 ns
- diffusion constant (nm2/ns): 1000.0
- emission of a delta amplitude 1000 molecules, time 0ns
- distance tx-rx 1µm





- simulation time 5000 μs Sampling Period 1000 ns
- diffusion constant (nm2/ns): 1000.0
- emission of a delta amplitude 1000 molecules, time Ons
- distance tx-rx 1µm





B)