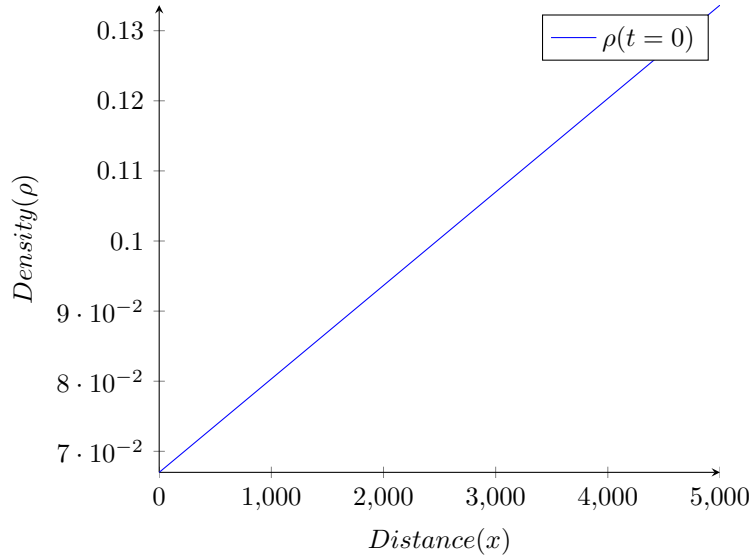


# LWR Equation Simulation

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## 1 Model presentation

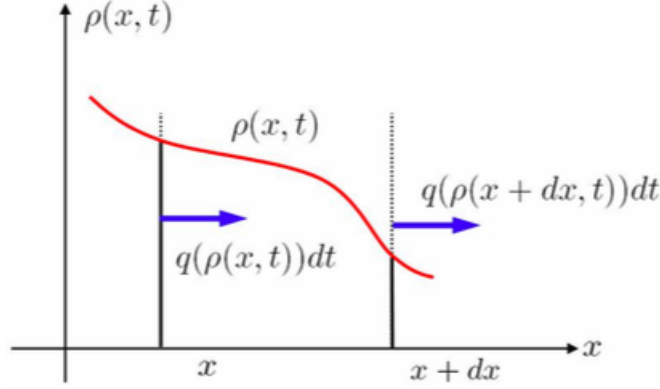


**Figure 1:** The vehicle density road system, typically the initial state

Our system expresses the vehicle density(number of vehicles per meter) on a 5000-meter-long road. This quantity varies with space( $x$ ) and time( $t$ ), and we use  $\rho(x, t)$  to denote it. In this case, we use the *Lighthill – Whitham – Richards(LWR)* PDE to study the system. To observe the change of vehicle density on different spaces with time going by, we set the initial density(when  $t=0$ ) of the road as shown in the Figure 1.

In order to quantify the evolution of the density of vehicles on the road, we use a mass balance for a small control volume of length  $dx$  in the road. Following Figure 2, we have four terms in the balance:

- 1)  $\rho(x, t)dx$  number of vehicles in the control volume  $[x, x + dx]$  at  $t$
- 2)  $\rho(x, t + dt)dx$  number of vehicles in the control volume  $[x, x + dx]$  at  $t + dt$
- 3)  $q(\rho(x, t))dt$  number of vehicles entering the control volume  $[x, x + dx]$  between  $t$  and  $t + dt$  through  $x$
- 4)  $q(\rho(x + dx, t))dt$  number of vehicles entering the control volume  $[x, x + dx]$  between  $t$  and  $t + dt$  through  $x + dx$



**Figure 2:** Illustration of the mass balance for the control volume  $[x, x + dx]$

Equating the four terms in the balance, we obtain:

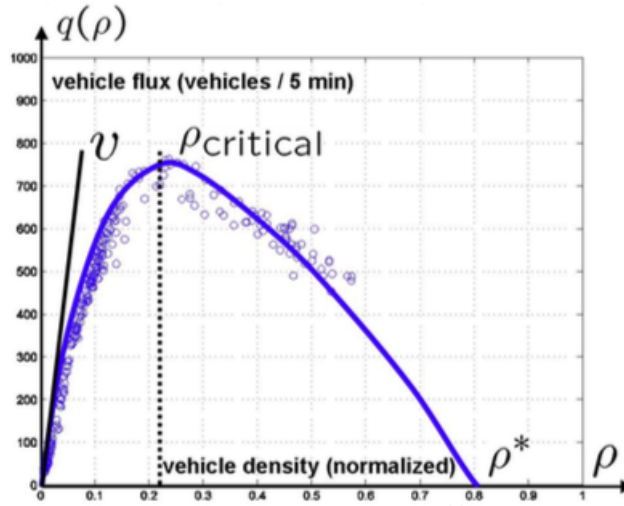
$$(\rho(x, t + dt) - \rho(x, t))dx = (q(\rho(x, t)) - q(\rho(x + dx, t)))dt \quad (1)$$

Dividing by  $dt$  and  $dx$ , and taking the limit as  $dt \rightarrow 0$  and  $dx \rightarrow 0$ , we obtain:

$$\frac{\partial \rho(x, t)}{\partial t} + \frac{\partial (q(\rho(x, t)))}{\partial x} = 0 \quad (2)$$

This PDE can alternatively be rewritten as:

$$\frac{\partial \rho(x, t)}{\partial t} + q'(\rho(x, t)) \frac{\partial (\rho(x, t))}{\partial x} = 0 \quad (3)$$



**Figure 3:** Greenshield model

*Greenshield Model* is an empirical measurement of the phenomenological law  $q$  with the density  $\rho$  as shown in figure 3. Each of the dots is one measurement. The solid curve is a fit of the

measurement. As can be seen, for small vehicle densities, the flux function increases almost linearly with the density (slope  $v$ ). It reaches a maximum for a critical density, called  $\rho_{critical}$ . For higher densities, it decreases until it finally reaches zero for a density  $\rho^*$  called jam density.

According to figure 3, *Greenshield flux function* is given by:

$$q(\rho) = v\rho(1 - \frac{\rho}{\rho^*}) \quad (4)$$

where  $\rho^*$  is the *jam density* and  $v$  is the *free flow velocity*.

Combine equation (3) and (4), we get the final equation for our modeling:

$$\frac{\partial \rho(x, t)}{\partial t} + v(1 - \frac{2\rho(x, t)}{\rho^*}) \frac{\partial \rho(x, t)}{\partial x} = 0 \quad (5)$$

## 2 Implementation

```

1 import numpy as np
2 import matplotlib.pyplot as plt
3 from mpl_toolkits.mplot3d import Axes3D
4
5 """
6 Data structure creation and initialisation
7 """
8 # Global parameters
9 T_final=3600 # time duration
10 N_t=500 # number of time steps: the time step value dt is computed later
11 X_final=5000 # road length
12 N_x=100 # number of space steps: the space step value dx is computed later
13 rho0=0.2 #jam density in Greenshield flux function(number of vehicles/m)
14 v0=15 #free flow velocity(m/s)
15
16 # structures for visualization and computation of time/space steps
17 T,dt=np.linspace(0,T_final,num=N_t,endpoint=True,retstep=True)
18 X,dx=np.linspace(0,X_final,num=N_x,endpoint=True,retstep=True)
19 #set the range and sample number of time(T) [0,3600]seconds in every 7.2 seconds
20 #set the range and sample number of distance(X) [0,5000]meters in every 50 meters
21 print("dx = ",dx," dt = ",dt)
22
23 # structure for simulation: density as a function of space and time
24 rho = np.zeros((N_x,N_t))
25 # initialization of the density at time 0 with a continuous function
26 for x in range(int(N_x/2)):
27     rho[x][0] = rho0/3+rho0*x/N_x/3 # from rho0/3 to rho0/2
28 for x in range(int(N_x/2),N_x):
29     # rho[x][0] = rho0/3+rho0*x/N_x/3 # from rho0/2 to rho0*2/3
30     rho[x][0] = rho0/3+rho0*x/N_x/3
31 print("t = 0")
32 plt.plot(X,rho[:,0])
33 plt.show()#plot the density in the range of x at t=0
34
35
36 """

```

```

37 Start the main simulation loop
38 Note that the naive Euler integration scheme is ALWAYS numerically unstable
39 Learn about Von Neumann stability analysis
40 And use the simple (stable) Lax scheme
41 But stability needs the Courant Friedrichs Levy condition to be verified
42 Trick: if unstable, decrease value of dt
43 """
44 for t in range(N_t-1): # at timestep t, compute rho at t+1
45     for x in range(1,N_x-1):
46         #calculate density at t+1 on i based on density at t on i and i+1
47         dr = v0*dt/dx*(2*rho[x][t]/rho0-1)*(rho[x+1][t]-rho[x-1][t])/2
48         r = (rho[x-1][t]+rho[x+1][t])/2 + dr #this is Lax Scheme
49         rho[x][t+1] = r
50     # for x==N_x, we take the derivative backward
51     x = 0
52     dr = v0*dt/dx*(2*rho[x][t]/rho0-1)*(rho[x+1][t]-rho[x][t])
53     r = (rho[x][t]+rho[x+1][t])/2 + dr
54     rho[x][t+1] = r
55     x = N_x-1
56     dr = v0*dt/dx*(2*rho[x][t]/rho0-1)*(rho[x][t]-rho[x-1][t])
57     r = (rho[x][t]+rho[x-1][t])/2 + dr
58     rho[x][t+1] = r
59     if ((t+1)%int(N_t/5))==int(N_t/5)-1:
60         print("t = ",7.2*t)
61         plt.plot(X,rho[:,t])
62         plt.show()#plot the density in the range of x at t
63         =705.6,1425.6,2145.6,2865.6,3585.6
64
65 X,T = np.meshgrid(X,T)
66 Z = rho.reshape(X.shape)
67
68 fig=plt.figure()
69 ax=Axes3D(fig)
70 ax.plot_surface(X,T,Z, rstride=1, cstride=1, cmap='rainbow')
71 plt.show()#plot the 3d figure

```

Listing 1: Python Code

The most important part to determine stability of the system lies in line 47 and 48 in the Listing 1. A stable system would have the same code in Listing 1 which is:

```

1 dr = v0*dt/dx*(2*rho[x][t]/rho0-1)*(rho[x+1][t]-rho[x-1][t])/2
2 r = (rho[x-1][t]+rho[x+1][t])/2 + dr

```

Listing 2: Stable Code

Whereas an unstable system would have the below code to replace the stable one:

```

1 dr = v0*dt/dx*(2*rho[x][t]/rho0-1)*(rho[x+1][t]-rho[x][t])
2 r = rho[x][t] + dr

```

Listing 3: Unstable Code

The stable one uses *Lax – Friedrichs* method, a numerical method for the solution of hyperbolic partial differential equations based on finite differences. Whereas the unstable one uses *Euler* method, a first order numerical procedure for solving ordinary differential equations with a given initial value.

To understand the difference between these two methods intuitively, we will take the *AdvectionEquation* as an example:

$$\frac{\partial u}{\partial t} = -v \frac{\partial u}{\partial x} \quad (6)$$

To express equation(6) with *Euler* method, we will get:

$$\frac{u_x^{t+1} - u_x^t}{\Delta t} = -v \left( \frac{u_{x+1}^t - u_x^t}{\Delta x} \right) \quad (7)$$

Whereas to express equation(6) with *Lax – Friedrichs* method, we will get:

$$\frac{u_x^{t+1} - \frac{1}{2}(u_{x+1}^t - u_{x-1}^t)}{\Delta t} = -v \left( \frac{u_{x+1}^t - u_{x-1}^t}{2\Delta x} \right) \quad (8)$$

We can use *Von Neumann* stability analysis to check if a numerical scheme is stable without computation. In numerical analysis, *Von Neumann* is a procedure used to check the stability of finite difference schemes as applied to linear partial differential equations. The analysis is based on the *Fourier* decomposition of numerical error. The stability of numerical schemes is closely associated with numerical error. A finite difference scheme is stable if the errors made at one time step of the calculation do not cause the errors to be magnified as the computations are continued. A neutrally stable scheme is one in which errors remain constant as the computations are carried forward. If the errors decay and eventually damp out, the numerical scheme is said to be stable. If, on the contrary, the errors grow with time the numerical scheme is said to be unstable. *Von Neumann* analysis is often used in place of a more detailed stability analysis to provide a good guess at the restrictions (if any) on the step sizes used in the scheme because of its relative simplicity.

In *Von Neumann* stability analysis, we assume:

$$u_x^t = \xi^t e^{ikx\Delta x} \quad (9)$$

$\xi$  is the amplification factor and  $k$  is wave number.  $\xi = \xi(k)$ . A numerical scheme is unstable if  $|\xi(k)| > 1$  and is stable if  $|\xi(k)| \leq 1$ .

First we investigate *Eular* method, from equation(7) we get:

$$u_x^{t+1} = u_x^t + \frac{v\Delta t}{\Delta x} (u_{x+1}^t - u_x^t) \quad (10)$$

Put equation(9) into (10), we get:

$$\xi = 1 + \frac{v\Delta t}{\Delta x} (\exp(ik\Delta x) - 1) \Rightarrow \xi = 1 + \frac{v\Delta t}{\Delta x} (\cos(ik\Delta x) + i\sin(ik\Delta x) - 1) \quad (11)$$

We get the amplitude  $|\xi| = \sqrt{\xi\xi^*}$ , where  $\xi^*$  is the conjugate complex of  $\xi$ :

$$|\xi| = \sqrt{[(\cos(ik\Delta x) - 1) \frac{v\Delta t}{\Delta x} + 1]^2 + \sin^2(ik\Delta x)} > 1 \quad (12)$$

which means *Eular* method is unconditionally unstable.

Then we investigate *Lax – Friedrichs* method in the same way, from equation(8) we get:

$$u_x^{t+1} = \frac{1}{2}(u_{x+1}^t - u_{x-1}^t) - \frac{v\Delta t}{2\Delta x} (u_{x+1}^t - u_{x-1}^t) \quad (13)$$

Put equation(9) into (13), we get:

$$\xi = \frac{1}{2}(\exp(ik\Delta x) + \exp(-ik\Delta x)) + \frac{v\Delta t}{2\Delta x}(\exp(ik\Delta x) - \exp(-ik\Delta x)) \Rightarrow \xi = \cos(k\Delta x) - i\frac{v\Delta t}{\Delta x}\sin(k\Delta x) \quad (14)$$

We get the amplitude:

$$|\xi| = \sqrt{\cos^2(k\Delta x) + \left(\frac{v\Delta t}{\Delta x}\right)^2 \sin^2(k\Delta x)} \quad (15)$$

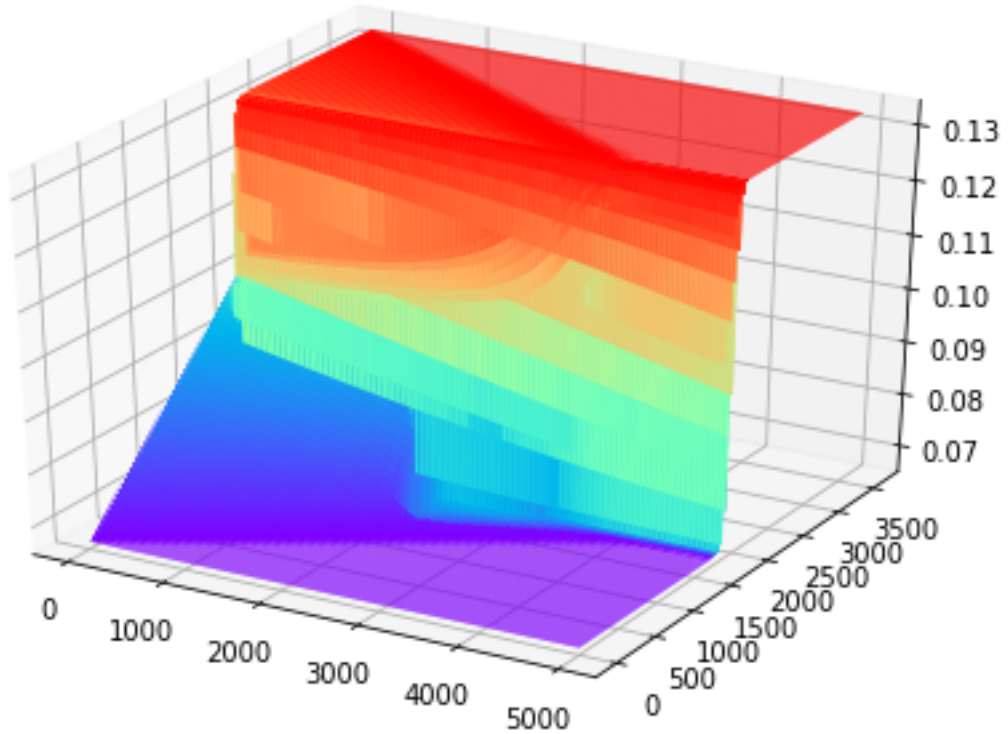
which means *Lax – Friedrichs* method is conditionally stable if:

$$\frac{|v|\Delta t}{\Delta x} \leq 1 \quad (16)$$

Equation(16) is called *Courant Friedrichs Levy* condition. It is a famous stability condition in numerical mathematics and is valid for many physical applications, also in inhomogenous nonlinear cases like Hydrodynamics (with v as sound speed), MHD (with v as Alfven velocity), etc.

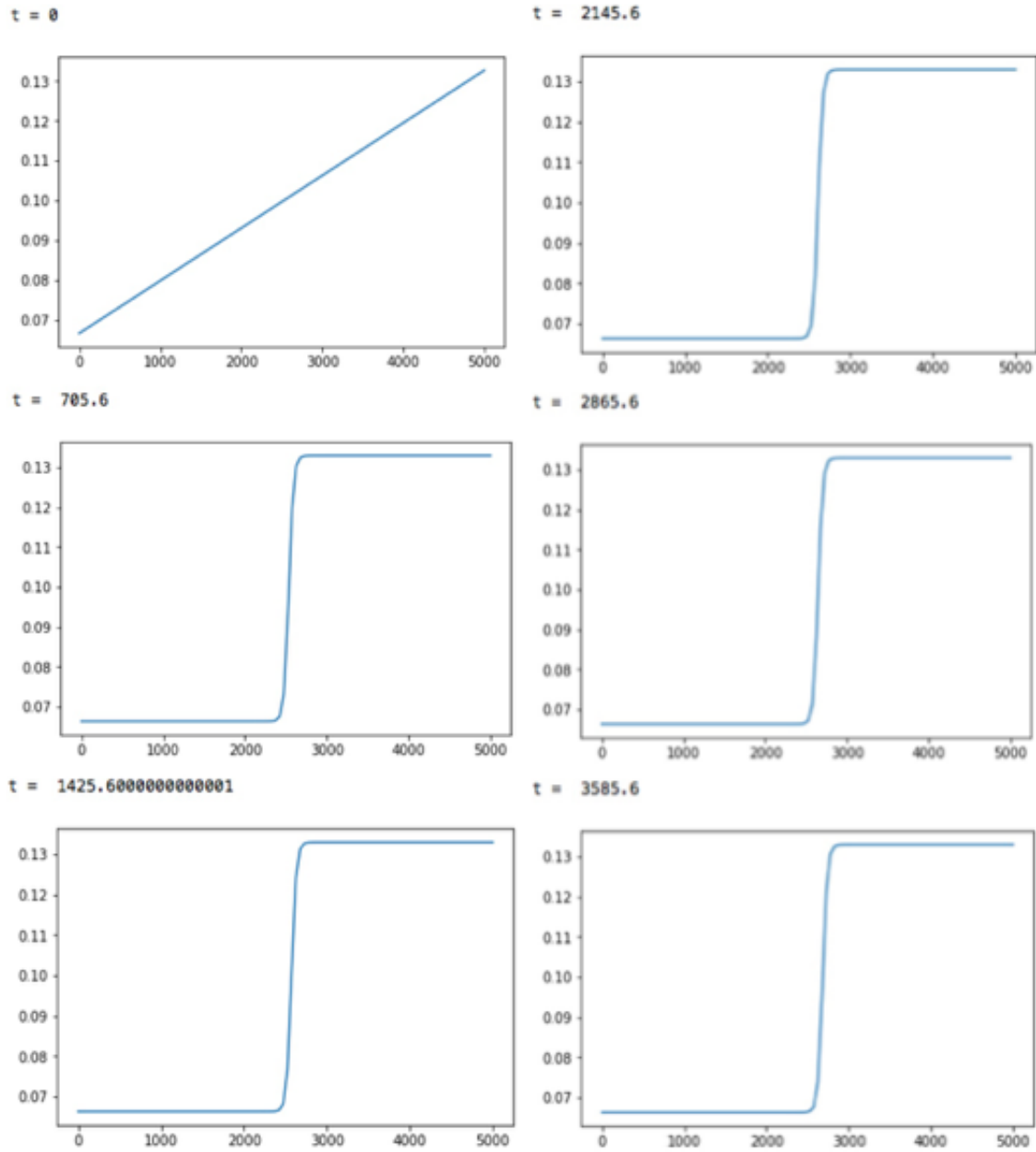
### 3 Results

For the stable code, we get final plot like this:



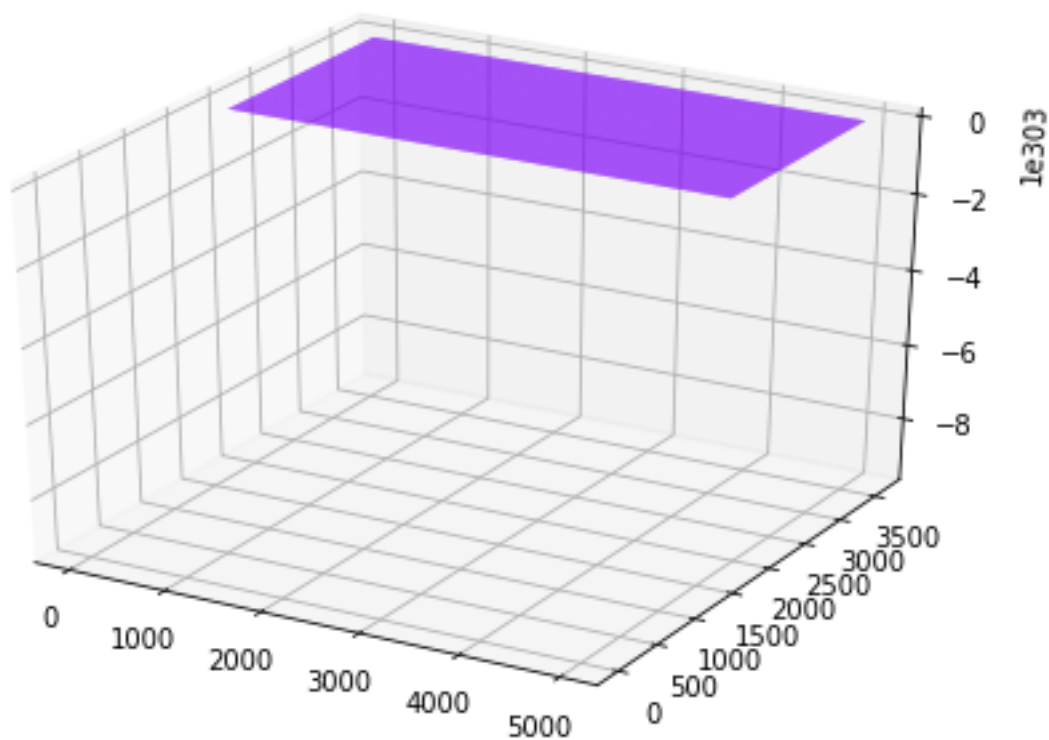
**Figure 4:** Final plot for stable code

And the density plot in the range of  $x$  at  $t=705.6, 1425.6, 2145.6, 2865.6, 3585.6$  respectively:



**Figure 5:** Density plot in the range of  $x$  for stable code

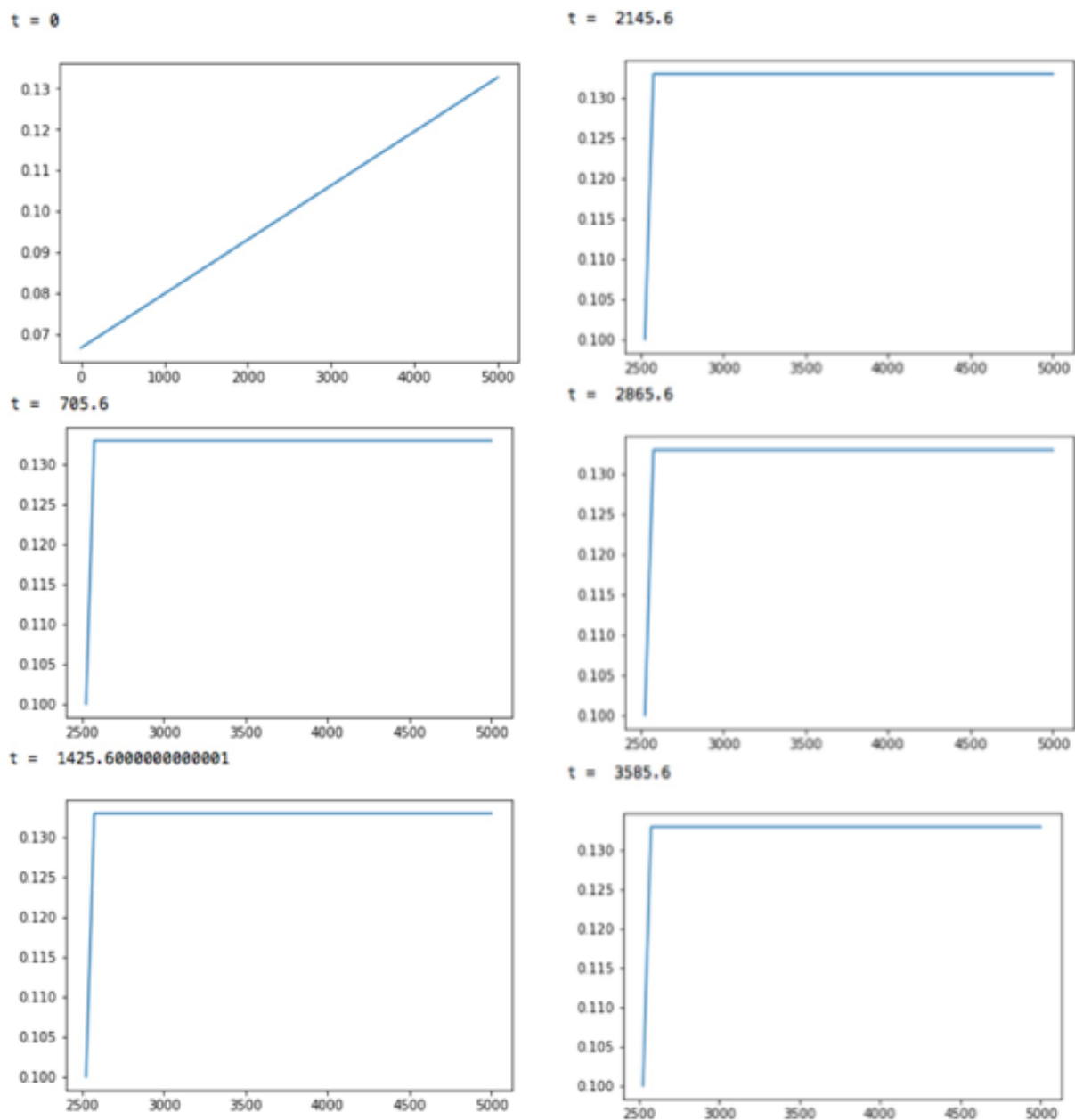
For the unstable code, we get final plot like this:



**Figure 6:** Final plot for unstable code

And the density plot in the range of  $x$  at  $t=705.6, 1425.6, 2145.6, 2865.6, 3585.6$  respectively:





**Figure 7:** Density plot in the range of  $x$  for unstable code

## 4 Interpretation

As we can see from the Result part, we can get reasonable outputs from the stable code using *Lax – Friedrichs* method. We understand that based on *LWR PDE* and *Greenshield Model*, we will get a shock wave in the middle of the road if we start with linearly increasing density on the road at time zero.

However, when it comes to the unstable code using *Eular* method, we get false outputs which make no sense. The error in each step tends to accumulate to infinite, and make the density far beyond the reasonable range.

## 5 Conclusion

From this project, we learn the *LWR PDE* and *Greenshield Model* and understand the physical meaning behind them. We learn how to describe a system and present a model. We also learn how to implement a model and do simulation with Python. And most importantly, we have realized the significance of numerical stability when doing such kind of modeling.

With the development of the code with *Lax – Friedrichs* method instead of *Eular* method, everything is aligned well in theory and practice.

As for the difficulty, it would be not realizing the numerical stability issue at the very beginning. Once the problem was solved, everything went out smoothly.