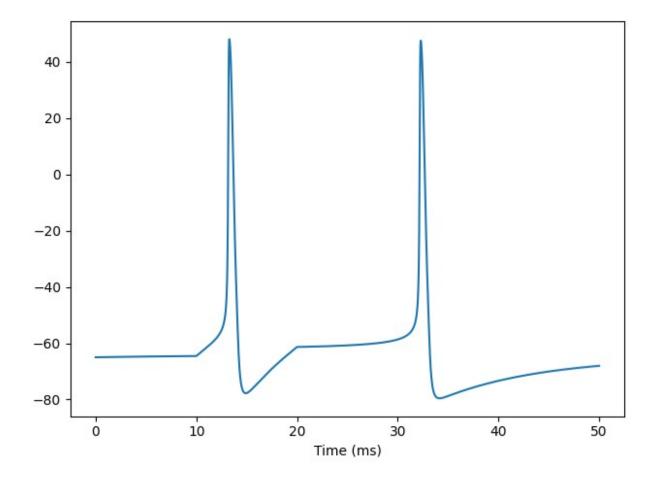
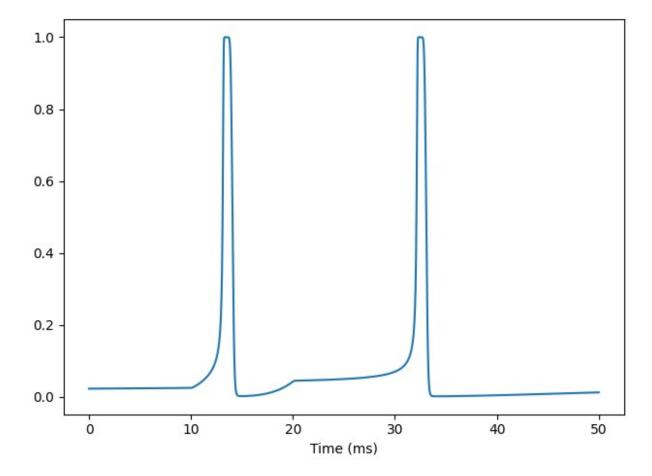
```
import numpy as np
import matplotlib.pyplot as plt
from types import SimpleNamespace
def euler integrate(
    derivs,
    χ0,
    t,
):
    x = np.empty((len(t), len(x0)))
    x[0] = x0
    for k in range(len(t) - 1):
        dt = t[k + 1] - t[k]
        x[k + 1] = x[k] + dt * derivs(t[k], x[k])
    return x
def RC derivative(tau, I):
    """f(x, t)"""
    def deriv(t, x):
        dx = -1 / tau * x + I(t)
        return np.array([dx])
    return deriv
def plot trajectory(
    t: np.ndarray, V: np.ndarray, ylab="", xlab="Time (ms)", title:
str = ""
):
    plt.figure()
    plt.plot(t, V)
    plt.xlabel(xlab)
    plt.ylabel(ylab)
    plt.title(title)
    plt.tight layout()
def alpha m(V):
    # Avoid division by zero
    denominator = 1 - np.exp(-(V + 54) / 4)
    return 0.32 * (V + 54) / denominator
def beta m(V):
    # Avoid division by zero
```

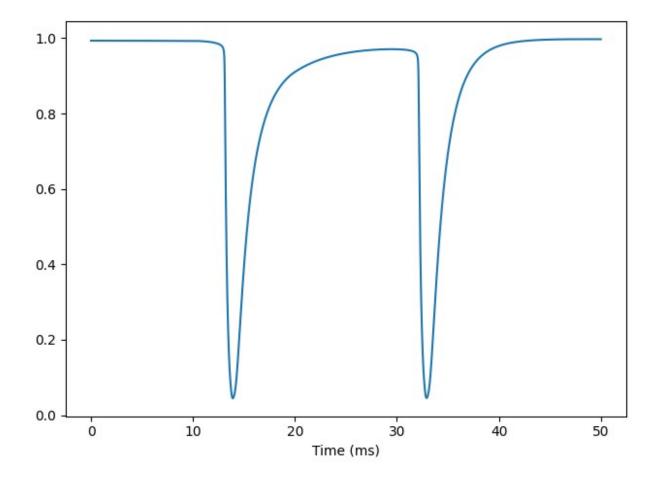
```
denominator = np.exp((V + 27) / 5) - 1
    return 0.28 * (V + 27) / denominator
def alpha h(V):
    # Prevent overflow in exp
    V_{clipped} = np.clip(-(V + 50) / 18, -500, 500)
    return 0.128 * np.exp(V clipped)
def beta h(V):
    # Prevent overflow in exp
    V \text{ clipped} = \text{np.clip}(-(V + 27) / 5, -500, 500)
    return 4 / (1 + np.exp(V clipped))
def alpha n(V):
    # Avoid division by zero
    denominator = 1 - np.exp(-(V + 52) / 5)
    return 0.032 * (V + 52) / denominator
def beta n(V):
    # Prevent overflow in exp
    V \text{ clipped} = \text{np.clip}(-(V + 57) / 40, -500, 500)
    return 0.5 * np.exp(V clipped)
def tau x(V, alpha x, beta x):
    # Avoid division by zero
    denominator = alpha x(V) + beta_x(V)
    return 1 / denominator
def x_inf(V, alpha_x, beta_x):
    # Avoid division by zero
    denominator = alpha x(V) + beta x(V)
    return alpha x(V) / denominator
def x deriv(V, x, alpha x, beta x):
    return -1 / tau x(V, alpha x, beta x) * (x - x inf(V, alpha x,
beta x))
# m: Sodium activation gate
# h: Sodium inactivation gate
def I NA(cfg, m, h, V):
    return cfg.gNa * m**3 * h * (V - cfg.ENa)
```

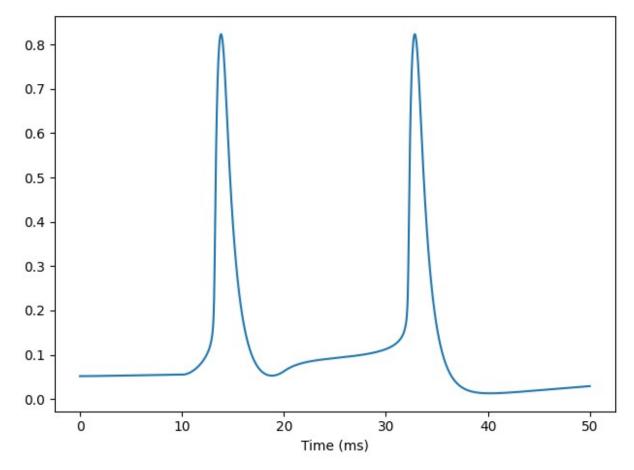
```
# n: Potassium activation gate
def I K(cfg, n, V):
    return cfg.gK * n**4 * (V - cfg.EK)
def I L(cfg, V):
    return cfg.gL * (V - cfg.EL)
def C m deriv(cfg, I ext, V, m, h, n):
    return I ext - I NA(cfg, m, h, V) - I K(cfg, n, V) - I L(cfg, V)
def hh derivatives(cfg, I):
    def derivs(t, x):
        V, m, h, n = x \# x = [V, m, h, n]
        # Calculate the time constants for the gating variables
        tau m = tau x(V, alpha m, beta m)
        tau h = tau x(V, alpha h, beta h)
        tau n = tau \times (V, alpha n, beta n)
        # Calculate the steady-state values for the gating variables
        m inf = x inf(V, alpha m, beta m)
        h inf = x inf(V, alpha h, beta h)
        n inf = x inf(V, alpha n, beta n)
        # derivative of the gating variables
        m_{deriv} = (m_{inf} - m) / tau_{m}
        h_deriv = (h_inf - h) / tau_h
        n deriv = (n inf - n) / tau n
        # Calculate the ionic currents
        g na = cfg.gNa * m**3 * h
        g k = cfg.gK * n**4
        g l = cfg.gL
        I Na = g na * (V - cfg.ENa)
        I K = g k * (V - cfg.EK)
        I_L = g_l * (V - cfg.EL)
        v deriv = (I(t) - I Na - I K - I L) / cfg.Cm
        return np.array([v deriv, m deriv, h deriv, n deriv])
    return derivs
def simulate hh(cfg, I, T=200.0, dt=0.025, v0=-65.0):
    t = np.arange(0.0, T + dt, dt)
```

```
m_inf = x_inf(v0, alpha_m, beta_m)
    h inf = x inf(v0, alpha h, beta h)
    n_inf = x_inf(v0, alpha_n, beta_n)
    x0 = np.array([v0, m inf, h inf, n inf])
    traj = euler integrate(hh derivatives(cfg, I), x0, t)
    return {"t": t, "V": traj[:, 0], "m": traj[:, 1], "h": traj[:, 2],
"n": traj[:, 3]}
cfg = SimpleNamespace(**{})
cfq.dt = 0.1
cfg.T = 50
cfg.t = np.arange(0.0, cfg.T + cfg.dt, cfg.dt)
cfg.tau = 20
# These variables are the parameters for the Hodgkin-Huxley model
cfg.Cm = 1.0 # Membrane capacitance
cfg.gNa = 50.0 # Sodium conductance
cfg.gK = 10.0 # Potassium conductance
cfg.gL = 0.1 # Leak conductance
cfg.ENa = 50.0 # Sodium reversal potential
cfg.EK = -90.0 # Potassium reversal potential
cfg.EL = -65.0 # Leak reversal potential
# Initial conditions
cfg.V 0 = -65.0
cfg.m_0 = x_inf(cfg.V_0, alpha_m, beta_m)
cfg.h 0 = x inf(cfg.V 0, alpha h, beta h)
cfg.n 0 = x inf(cfg.V 0, alpha n, beta n)
cfg.x_0 = [cfg.V_0, cfg.m_0, cfg.h_0, cfg.n 0]
T = 50.0
I amp = 3.0
I = lambda t: I amp if 10.0 \ll t < 20.0 else 0.0
data = simulate hh(cfg, I, T)
plot_trajectory(data["t"], data["V"])
plot_trajectory(data["t"], data["m"])
plot_trajectory(data["t"], data["h"])
plot trajectory(data["t"], data["n"])
```









```
v_array = np.linspace(-100, 40, 100)

tau_m_array = tau_x(v_array, alpha_m, beta_m)
tau_n_array = tau_x(v_array, alpha_n, beta_n)
tau_h_array = tau_x(v_array, alpha_h, beta_h)

x_inf_m_array = x_inf(v_array, alpha_m, beta_m)
x_inf_n_array = x_inf(v_array, alpha_n, beta_n)
x_inf_h_array = x_inf(v_array, alpha_h, beta_h)

plot_trajectory(v_array, tau_m_array, xlab="Voltage (mV)", ylab="Tau")
plot_trajectory(v_array, tau_n_array, xlab="Voltage (mV)", ylab="Tau")
plot_trajectory(v_array, tau_h_array, xlab="Voltage (mV)", ylab="Tau")
```

