

Adaptive Exponential Rosenbrock–Euler Method (ExpRESS–Euler)’s report

The ODE System

$$dv/dt = k(v - v_r)(v - v_t) - w + I_n$$

$$dw/dt = a[b(v - v_r) - w]$$

Parameters

C : Membrane capacitance (value: $C = 100$, units: typically pF or scaled).

v : Neuron membrane potential (in millivolts, mV).

w : Recovery current (in picoamperes, pA , or scaled units).

t : Time (in milliseconds, ms).

v_r : Resting membrane potential (value: $v_r = -60\text{mV}$).

k : Scaling factor for the quadratic term (value: $k = 0.7$).

v_t : Threshold membrane potential (value: $v_t = -40\text{mV}$).

I_n : Input current (values: $I_n = 0, \text{for } t < 101\text{ms}, I_n = 70, \text{for } t \geq 101\text{ms}$, units: pA).

a : Time scale of recovery (value: $a = 0.03$, units: ms^{-1}).

b : Sensitivity of recovery to membrane potential (value: $b = -2$, dimensionless or scaled).

Initial Conditions

$$v(t = 0) = v_0, w(t = 0) = w_0$$

v_0 and w_0 constants based on the state of the neuron.

Reset Mechanism

consider a membrane potential threshold $v_{\text{peak}} = 35 \text{ mV}$, when the membrane potential reaches v_{peak} :

- we set $v(t) = -50$ mV
 - set $w(t) = w(t) + 100$
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Adaptive Exponential Rosenbrock–Euler Method

The Adaptive Exponential Rosenbrock-Euler method (ExpRESS-Euler) is a numerical technique for solving ordinary differential equations (ODEs), particularly those that are stiff, meaning they contain vastly different timescales. It combines the concepts of exponential integrators, Rosenbrock methods, and adaptive step-size control.

Method's Formula

In standard ODE form:

$$dy/dt = f(t, y)$$

The **ExpRESS–Euler** updates the solution using:

$$y_{n+1} = y_n + h \ \varphi_1(h * J_n) [f(t_n, y_n)]$$

Where:

y_n : current solution vector (e.g. $[v, w]$ in neuron models).

h is the Adaptive Step Size. It adjusts the time step size based on local error estimation or stiffness indicators, This keeps the method stable and efficient across varying dynamics.

$J_n \approx \partial f / \partial y$ = Jacobian (or an approximation).

$\varphi_1(z) = (e^z - 1)/z$, a *matrix function* that comes from exponential integrators.

How to solve ODE using this Method

Step 1: Compute Current Values

At current time t_n , compute:

- y_n : your state vector
- $f_n = f(t_n, y_n)$

Step 2: Approximate or Compute the Jacobian

$J_n = \partial f / \partial y$ (at t_n and y_n).

- Can be exact (symbolically or numerically) or approximated via finite differences.

Step 3: Evaluate the Matrix Function φ_1

Now comes the core:

$$\varphi_1(h * J_n)$$

How do you do this?

- Most practical way: Use a matrix identity:
- $\varphi_1(hJ) \approx (I - hJ)^{-1}$, this is a *first-order* approximation using a resolvent (comes from the exponential series).

Step 4: Update the Solution

Use the main formula:

$$y_{n+1} = y_n + h \cdot \varphi_1(h * J_n) [f(t_n, y_n)]$$

or you can use substitute $\varphi_1(h * J_n)$ by the resolvent in step 3.

Step 5: Estimate Error (For Adaptivity)

To control the step size adaptively:

1. Compute the next point using one full step:
2. $y_1 = y_n + h \cdot \varphi_1(h J) \cdot f$
3. Then compute it again using **two half steps**:
 - Step 1: $y_{n+1/2}$
 - Step 2: y_{n+1}

Then:

$$\text{error} = \|y_1 - y_2\|$$

where y_1 is the y calculated using 1 step size and y_2 is the y calculated using 2 step sizes

Step 6: Adjust Step Size h

If $\text{error} < \text{tolerance}$, accept the step. Otherwise, reduce h and retry.

Adapt h using:

$$h_{\text{new}} = h \cdot (\text{tol}/\text{error})^{(0.5)}$$

Clamp it between h_{min} and h_{max} to avoid insanity.

Step 7: Repeat Until Final Time.

The Python Code

```
1  import numpy as np
2  import matplotlib.pyplot as plt
3  from numpy.linalg import inv
4  import pandas as pd
5  import time
6
7  start_time = time.time()
8  # Constants
9  C = 100
10 k = 0.7
11 vr = -60
12 vt = -40
13 a = 0.03
14 b = -2
15 v_peak = 35
16 c = -50
17 d = 100
18
19 # Simulation setup
20 T = 1000 # total time (ms)
21 h = 0.25 # initial step size
22 tol = 0.5 # tolerance for adaptive step
23 fac_min, fac_max = 0.1, 5
24 h_min, h_max = 0.01, 2.0
25
26 # Initial conditions
27 t_vals = [0]
28 v_vals = [vr]
29 w_vals = [0]
30 h_vals = [h]
31
32 # Table data
33 recorded_data = []
34
35 t = 0
36 v = vr
37 w = 0
38
```

```

39 while t < T:
40     In = 0 if t < 101 else 70
41
42     # Current state
43     y = np.array([v, w])
44
45     # Compute f(y)
46     dvdt = (1 / C) * (k * (v - vr) * (v - vt) - w + In)
47     dwdt = a * (b * (v - vr) - w)
48     f = np.array([dvdt, dwdt])
49
50     # Compute Jacobian
51     dv_dv = (1 / C) * k * ((v - vr) + (v - vt))
52     dv_dw = -1 / C
53     dw_dv = a * b
54     dw_dw = -a
55     J = np.array([
56         [dv_dv, dv_dw],
57         [dw_dv, dw_dw]
58     ])
59
60     # First-order ExpRESS-Euler step
61     I = np.eye(2)
62     phi1 = inv(I - h * J)
63     y1 = y + h * phi1 @ f
64
65     # Half step + half step for error estimate (2nd-order Richardson)
66     h_half = h / 2
67     phi_half = inv(I - h_half * J)
68     y_half = y + h_half * phi_half @ f
69
70     v_half, w_half = y_half
71     dvdt_half = (1 / C) * (k * (v_half - vr) * (v_half - vt) - w_half + In)
72     dwdt_half = a * (b * (v_half - vr) - w_half)
73     f_half = np.array([dvdt_half, dwdt_half])
74     y2 = y_half + h_half * phi_half @ f_half
75
76     # Error estimate
77     err = np.linalg.norm(y1 - y2)
78     err = max(err, 1e-10) # Avoid divide-by-zero

```

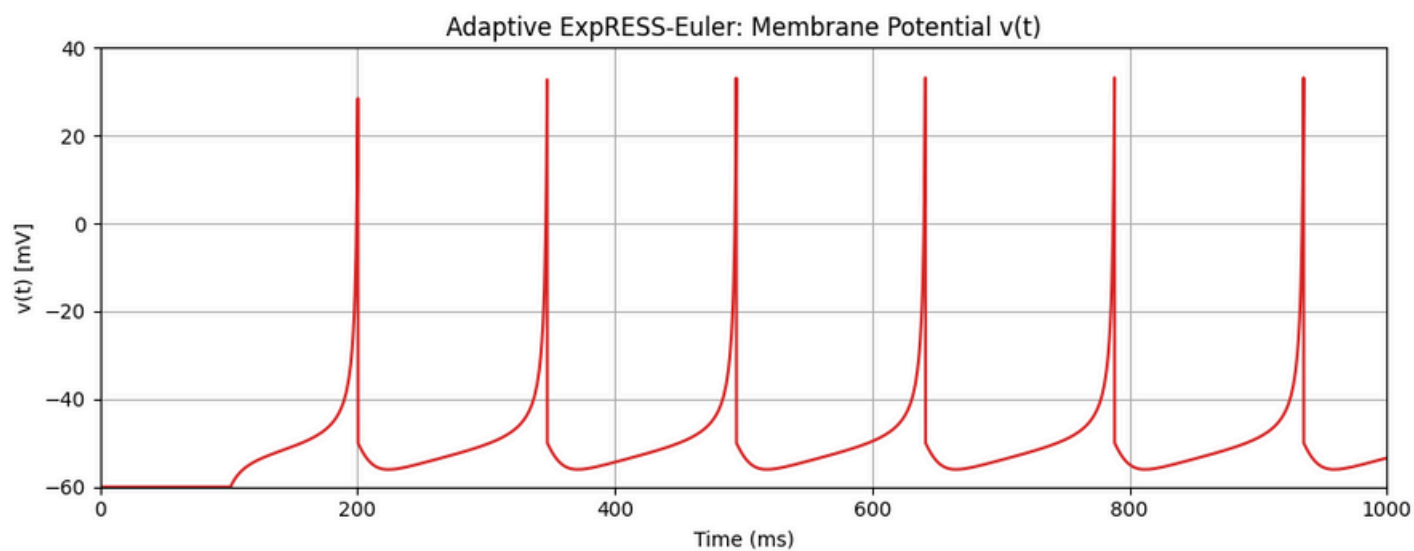
```

79
80     # Adaptive time step control
81     if err < tol:
82         t += h
83         v, w = y2 # accept higher-order result
84         if v >= v_peak:
85             v = c
86             w += d
87
88         t_vals.append(t)
89         v_vals.append(v)
90         w_vals.append(w)
91         h_vals.append(h)
92         recorded_data.append([t, v, w, h])
93
94         # Adjust h for next step
95         h = h * min(max((tol / err) ** 0.5, fac_min), fac_max)
96         h = min(max(h, h_min), h_max)
97     else:
98         # Reduce h and retry
99         h = h * max((tol / err) ** 0.5, fac_min)
100        h = max(h, h_min)
101
102    end_time = time.time()
103    # Print execution time
104    print(f"Execution time: {end_time - start_time:.8f} seconds")

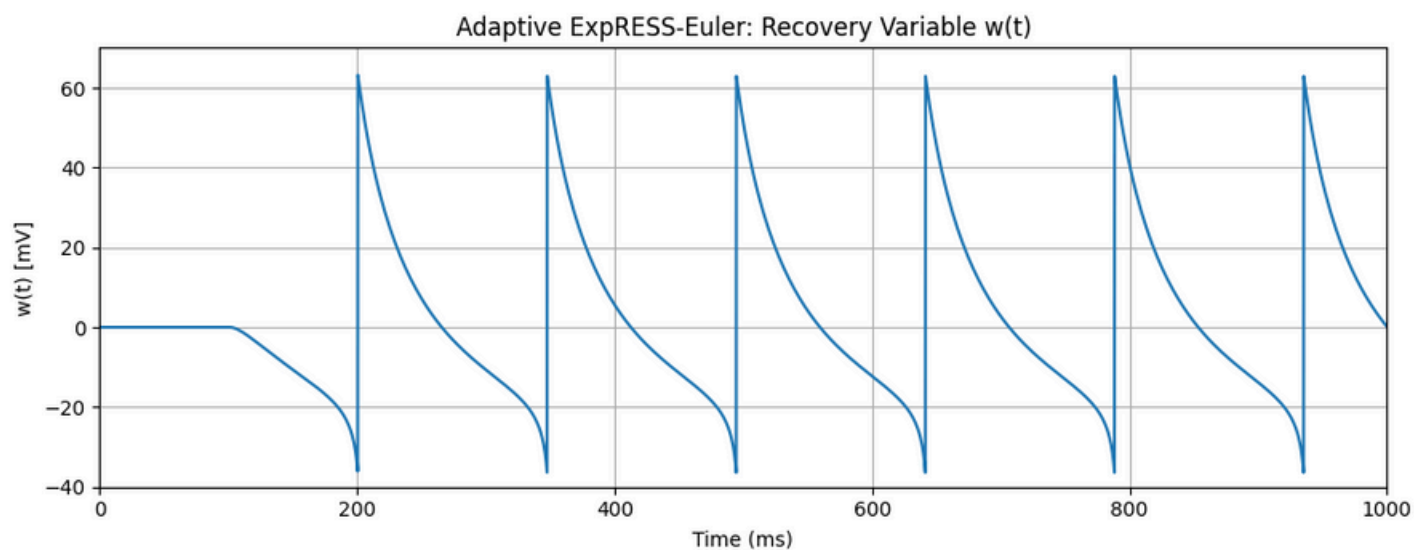
```

Code Output

V vs. t

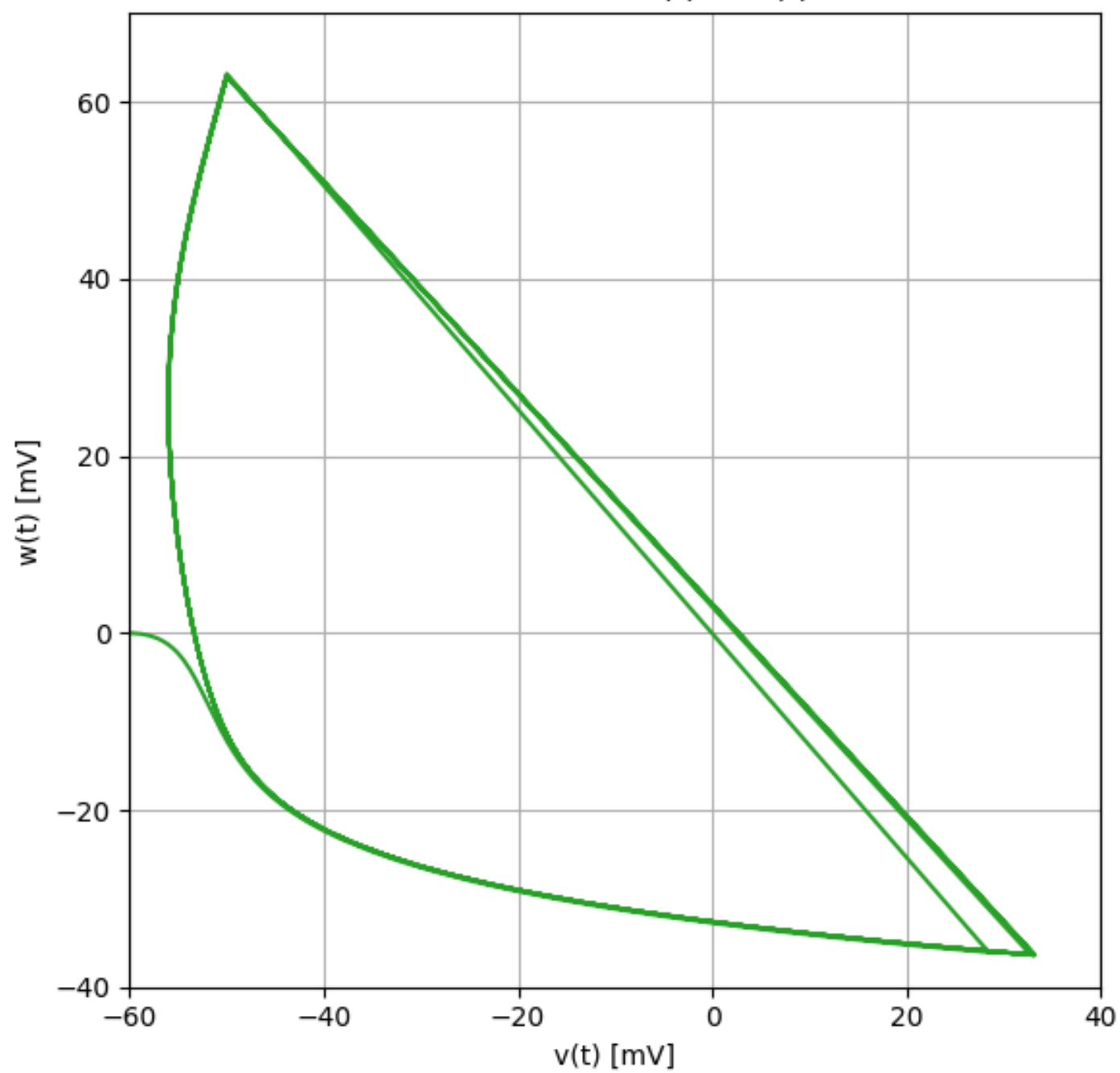


w vs. t



w vs. V

Phase Plane: $w(t)$ vs $v(t)$



	Time (ms)	v (mV)		w	Step size h
50	99.5	-60.000000	0.000000		2.0
51	101.5	-60.000000	0.000000		2.0
52	103.5	-58.843849	-0.102094		2.0
53	105.5	-57.917374	-0.310589		2.0
54	107.5	-57.156761	-0.598737		2.0
55	109.5	-56.519033	-0.946768		2.0
56	111.5	-55.974270	-1.339804		2.0
57	113.5	-55.501063	-1.766494		2.0
58	115.5	-55.083710	-2.218085		2.0
59	117.5	-54.710461	-2.687777		2.0
60	119.5	-54.372350	-3.170260		2.0

The error calculated relative to Explicit Euler Method

Time (ms)	v_sim	v_ref	v_err	v_err %	w_sim	w_ref	w_err	w_err %
0.0	-60.0000	-60.0000	0.0000	0.00%	0.000	0.0000	0.0000	—
250.0	-55.0123	-54.4819	0.5304	0.97%	5.950	6.2834	0.3334	5.31%
500.0	-51.0000	-50.6154	0.3846	0.76%	60.000	59.0910	0.9090	1.54%
750.0	-50.0000	-49.5530	0.4470	0.90%	-13.000	-12.4763	0.5237	4.20%
1000.0	-54.0000	-53.6973	0.3027	0.56%	2.000	1.5649	0.4351	27.80%

Pros and Cons of using this Method

Pros

- Handles Stiff Systems Well.
- Adaptive Time Stepping
- Stable at Larger Time Steps

Cons

- Requires Jacobian Computation, for high-dimensional systems, this becomes computationally expensive.
- Only First-Order Accurate
- Harder to Implement than Euler or RK, relatable $t_{bh}(T,T)$.

