# **Better Learning for ODEs**





## **Decomposing Sequences**

### We can address the first two issues using a reformulation

Let's consider the sequence of measurements  $\{\hat{y}_k\}_{k=0}^n$ 

- We can view it as a sequence of pairs  $\{(\hat{y}_{k-1}, \hat{y}_k)_{k=1}^n$
- lacksquare ... Each referring to a distinct ODE, i.e.  $\dot{y}_k = f(y_k, t, \omega)$
- lacksquare ...But all ODEs share the same parameter vector  $oldsymbol{\omega}$

### With this approach, we can reformulate the training problem as:

argmin<sub>$$\omega$$</sub>  $\sum_{k=1}^{n} L(y_k(\hat{t}), \hat{y}_k)$   
subject to:  $\dot{y}_k = f(y_k, t, \omega) \qquad \forall k = 1..n$   
 $y_k(\hat{t}_{k-1}) = \hat{y}_{k-1} \qquad \forall k = 1..n$ 



## **Decomposing Sequences**

### Let's examine again the new training problem:

argmin<sub>$$\omega$$</sub>  $\sum_{k=1}^{n} L(y_k(\hat{t}), \hat{y}_k)$   
subject to:  $\dot{y}_k = f(y_k, t, \omega)$   $\forall k = 1..n$   
 $y_k(\hat{t}_{k-1}) = \hat{y}_{k-1}$   $\forall k = 1..n$ 

There a few things to keep in mind:

- The approach is viable only if we have measurements for the full state
- ...And we are also assuming that the original loss is separable
- Finally, the new training problem is not exactly equivalent to the old one
- ...Since by re-starting at each step we are disregarding compound errors





## **Preparing the Data**

### Our implementation can naturally deal with the reformulation

We just need to properly prepare the data

■ Each ODE can be seen as a different example

```
In [2]: ns = len(data.index)-1
```

- The sequence for each example contains only two measurements
- ...Corresponding to consecutive evaluation points

```
In [3]: tr_T = np.vstack((data.index[:-1], data.index[1:])).T
    print(tr_T[:3])

[[0. 1.]
    [1. 2.]
    [2. 3.]]
```





## **Preparing the Data**

### Our implementation can naturally deal with the reformulation

We just need to properly prepare the data

■ The first measurement represents the initial state

■ The second to the final state, which we need to define a target tensor





[2.6543906 111

## **Training**

### Then we can perform training as usual

```
In [8]: %%time
    dRC = util.RCNablaLayer(tau_ref=10, vs_ref=10)
    euler = util.ODEEulerModel(dRC)
    history = util.train_ml_model(euler, [tr_y0, tr_T], tr_y, validation_split=0.0, epochs=400)
    util.plot_training_history(history, figsize=figsize)
```

200

epochs

250

300

350

400

Model loss: 0.0000 (training)

50

CPU times: user 1.21 s, sys: 245 ms, total: 1.46 s

100

150

Wall time: 1.02 s





0.00

## **Training**

### The results are the same as before (including estimation problems)

```
In [9]: print(f'tau: {tau:.2f} (real), {dRC.get_tau().numpy()[0]:.2f} (estimated)')
    print(f'Vs: {Vs:.2f} (real), {dRC.get_vs().numpy()[0]:.2f} (estimated)')

    tau: 8.00 (real), 8.50 (estimated)
    Vs: 12.00 (real), 12.00 (estimated)
```

### ...But there are significant computational advantages

Since we are using a shallow compute graph rather than a deep one

- The training time is much lower
- Potential vanishing/exploding gradient problems are absent

Since we now have multiple examples

- We can benefit from stochastic gradient descent
- We can use a validation set





## **Training**

### Just keep in mind that using a validation set will slow down the process

```
In [10]: %%time
         dRC = util.RCNablaLayer(tau ref=10, vs ref=10)
         euler = util.ODEEulerModel(dRC)
         history = util.train ml model(euler, [tr y0, tr T], tr y, validation split=0.2, epochs=400, pati
          util.plot training history(history, figsize=figsize)
           0.08
           0.06
           0.04
           0.02
           0.00
                           50
                                    100
                                              150
                                                                  250
                                                                             300
                                                                                       350
                                                                                                400
                                                        200
                                                        epochs
          Model loss: 0.0000 (training) 0.0000 (validation)
          CPU times: user 6.61 s, sys: 728 ms, total: 7.34 s
          Wall time: 5.81 s
```





## **Accuracy Issues**

### We are now ready to tackle our estimation issues

- lacktriangle We know we have trouble estimating the au parameter
- Intuitively, that should translate in trouble estimating the dynamic behavior

#### Let's check whether this is true

■ We prepare data structures to replicate our original run





## **Accuracy Issues**

### Then we can run Euler method directly using our model

As a side benefit, this will naturally use the estimate parameters

```
In [12]: run_y = euler.predict([run_y0, run_T])
```

Next, let's build a dataset with the original data and the predictions:

```
In [13]: data_euler = data.copy()
   data_euler['euler'] = run_y[0]
   data_euler.head()
```

#### Out[13]:

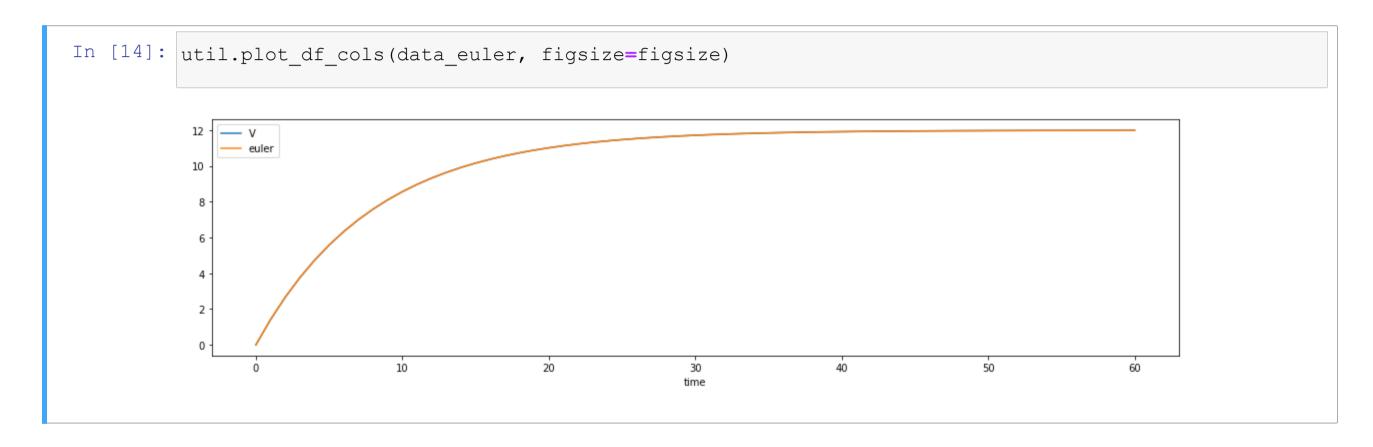
	V	euler
time		
0.0	0.000000	0.000000
1.0	1.410037	1.411475
2.0	2.654391	2.656847
3.0	3.752529	3.755661
4.0	4.721632	4.725165





## **Accuracy Issues**

### Finally, we can plot the two curves



We have a very good match!







## **Accuracy Issues?**

### We formulated the training problem in terms of curve fitting

- lacksquare I.e. we optimized au and  $V_s$  so as to obtain a close fitting curve
- ...Constructed using Euler method

### The problem is that Euler method is inaccurate

- If using wrong parameters will lead to a better fitting curve
- ...Our approach will not hesitate to do just that

### Is this a problem?

If we just care about the curve, not at all

■ It can actually be an advantage, if properly exploited

If we care about estimating parameters, then yes

■ ...But it also suggests an easy fix (using a more accurate integration method)





## **Improving Parameter Estimation**

### For sake of simplicity, we will keep using Euler method

...And we will just increase the number of steps to improve its accuracy

■ First, we introduce more evaluation points for each measurement pair

```
In [15]: nsteps = 11
    tr_T2 = np.vstack(np.linspace(data.index[:-1], data.index[1:], nsteps)).T
    print(tr_T2[:2])

[[0.    0.1    0.2    0.3    0.4    0.5    0.6    0.7    0.8    0.9    1. ]
        [1.    1.1    1.2    1.3    1.4    1.5    1.6    1.7    1.8    1.9    2. ]]
```

Second, we update the target sequences to match the size





## **Improving Parameter Estimation**

### Then, we can train as usual

```
In [17]: %%time
          dRC2 = util.RCNablaLayer(tau ref=10, vs ref=10)
          euler2 = util.ODEEulerModel(dRC2)
          history = util.train ml model(euler2, [tr y0, tr T2], tr y2, validation split=0.0, epochs=400)
          util.plot training history(history, figsize=figsize)
           0.06
           0.05
           0.04
           0.03
           0.02
           0.01
           0.00
                            50
                                      100
                                                 150
                                                                     250
                                                                                300
                                                                                           350
                                                                                                     400
                                                           200
                                                          epochs
```

Model loss: 0.0000 (training)

CPU times: user 2.32 s, sys: 292 ms, total: 2.61 s

Wall time: 1.94 s





## **Improving Parameter Estimation**

### This approach leads to considerably better estimates

```
In [18]: print(f'tau: {tau:.2f} (real), {dRC2.get_tau().numpy()[0]:.2f} (estimated)')
    print(f'Vs: {Vs:.2f} (real), {dRC2.get_vs().numpy()[0]:.2f} (estimated)')

    tau: 8.00 (real), 8.08 (estimated)
    Vs: 12.00 (real), 12.01 (estimated)
```

- The results can be improved by using additional steps
- ...Or by switching to a different integration method (e.g. RK4)

### Overall, when using this appraoch...

...It's important to be aware that integration methods are approximate

- This can easily lead to incorrectly estimated parameters
- Which may or may not be a problem, depending on your priorities



