

# **Comparison of different Data Driven Approaches for Li-Ion Battery Modeling**

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# Battery Modeling

- Li-Ion batteries are common in Electric vehicle battery packs nowadays
- Like many other dynamical systems, the electrochemical reactions are governed by multiple high dimensional differential equations
- This makes deriving a model of the battery a cumbersome task
- The advent of data-driven approaches for dynamical systems comes as a huge relief

# Goal



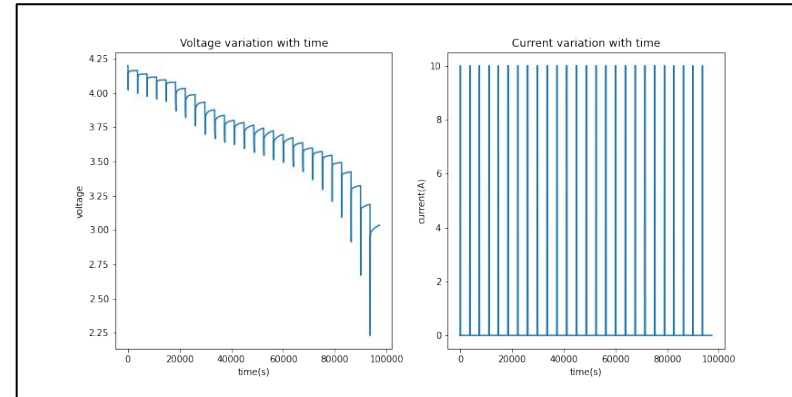
The goal of the project is model discovery of Li-Ion cells using both black and grey box modeling approaches

We are comparing and analysing

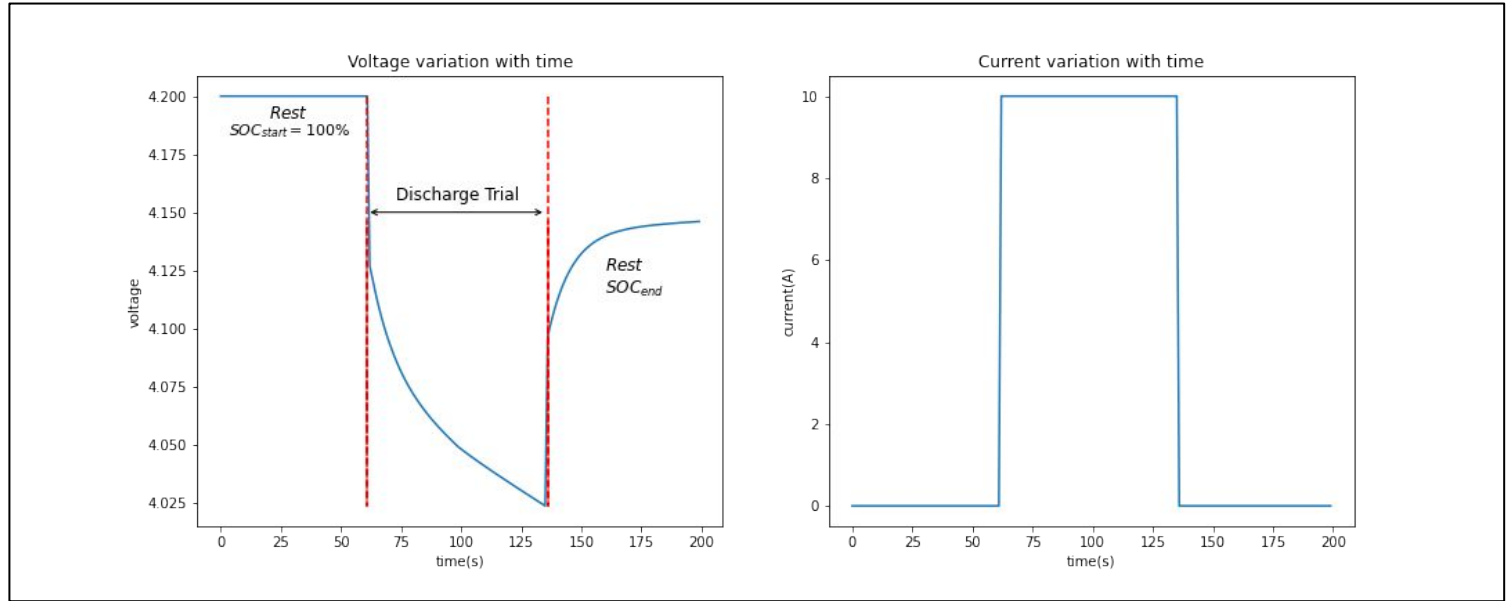
- SINDy - black box modeling
- Gaussian Process Regression - Grey box
- Genetic Algorithms - Grey box
- Regression - Grey box

# Pulse Discharge Experiment

- In this project, we are taking a data set of a Li-ion battery model which is obtained using a Pulse discharge experiment. The experiment involves applying a current discharge pulse for a certain period of time on a cell at regular intervals with a resting period in-between for measuring the cell voltage with a predefined sampling rate. The test is performed from a fully charged to completely drained battery state.



# Experiment



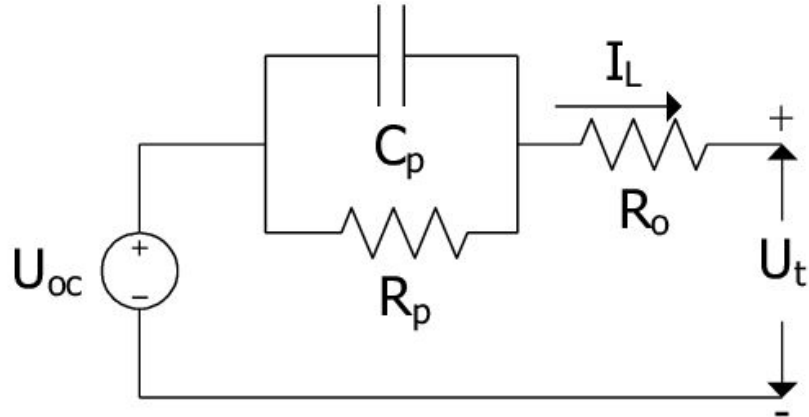
# Equivalent Circuit Modeling

The governing equation

$$U_t = U_{oc} - I_L R_0 - (I_L R_p (1 - e^{-t/(R_p C_p)}))$$

where,

- $U_{oc}$  = Open Circuit Voltage
- $R_0, R_p$  = Resistance
- $C_p$  = Capacitance
- $U_t$  = terminal voltage
- $I_L$  = input current



# GENETIC ALGORITHM



- > Genetic Algorithm (GA) is a search technique used in computing to find true or approximate solutions to optimization and search problems
- > Class of Evolutionary algorithms, that use techniques inspired from evolutionary biology
- > Developed to:
  - Understand the adaptive processes of natural systems
  - Design Artificial system to retain the robustness of natural system

# GENETIC ALGORITHM

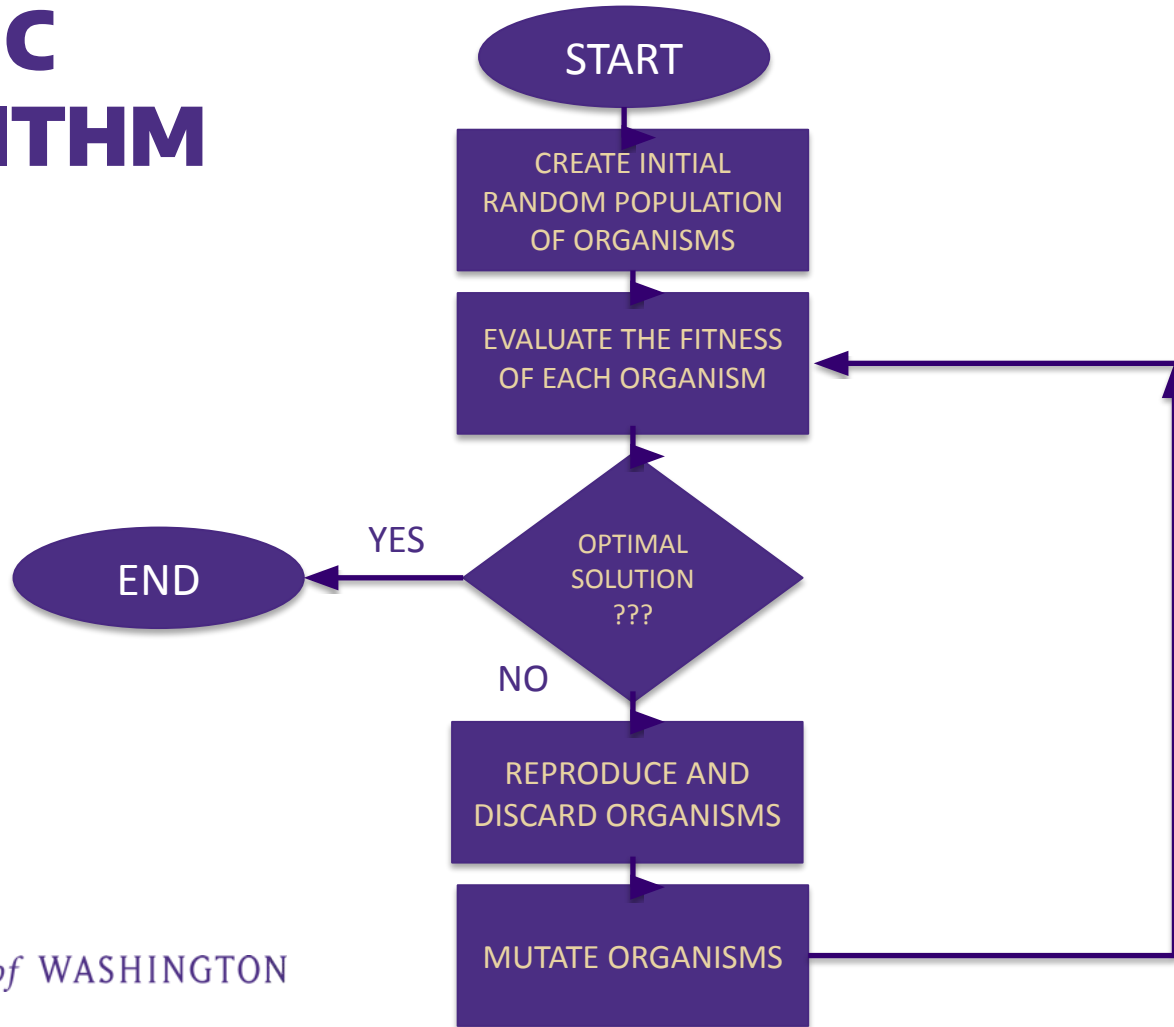


- > Basic idea is based on “Survival of the fittest”
- > Provides effective techniques for optimization and machine learning applications
- > Useful when search space is too large or complex
- > Examples in real world GA applications:
  - Antenna Design, Turbine engine design, Network Design, Data Mining, Control systems design



# GENETIC ALGORITHM

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# GENETIC ALGORITHM



```
def Genetic_Algorithm():
```

```
    initialize population
```

```
    evaluate the fitness of each individual in the population
```

```
    while optimal solution not found do:
```

```
        select fittest (best ranking) parents for reproduction
```

```
        perform mutation/crossover to breed new generation
```

```
        re-evaluate the fitness of the new population
```

```
    return (optimal solution)
```

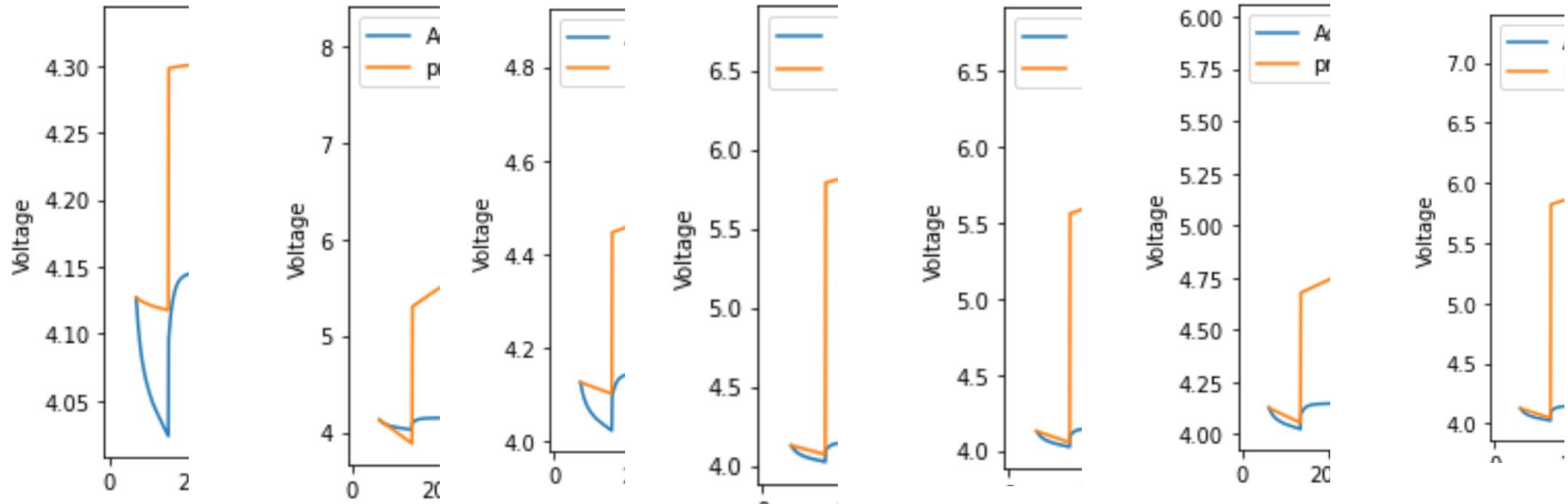
# GENETIC ALGORITHM

## > Battery Model

```
def battery(Em,R_0,R_1,C_1,R_2,C_2,R_3,C_3):  
    '''Equation of the model for the voltage is  
  
    
$$Em - I \cdot R_0 - (I \cdot R_1 (1 - \exp(-t/(R_1 \cdot C_1))))$$
  
    
$$+ (I \cdot R_2 (1 - \exp(-t/(R_2 \cdot C_2))))$$
  
    
$$+ (I \cdot R_3 (1 - \exp(-t/(R_3 \cdot C_3)))) - v\_data[t]$$
  
    ...  
    for t in range(62,1500):  
        if 61<t<136:  
            # for the time between t=62 and t=136 , the current I is 10 Amperes  
            return Em -10*R_0- ((10*R_1*(1- np.exp(-t/(R_1*C_1))))  
                                +(10*R_2*(1- np.exp(-t/(R_2*C_2))))  
                                +(10*R_3*(1- np.exp(-t/(R_3*C_3))))) - v_data[t]  
        else:  
            # for all other time instances , the current I is -10 Amperes  
            return Em +10*R_0- ((-10*R_1*(1- np.exp(-t/(R_1*C_1))))  
                                -(10*R_2*(1- np.exp(-t/(R_2*C_2))))  
                                -(10*R_3*(1- np.exp(-t/(R_3*C_3))))) - v_data[t]
```

# GENETIC ALGORITHM

Solutions of each generation



# GENETIC ALGORITHM

## Results

====gen0 best solutions====

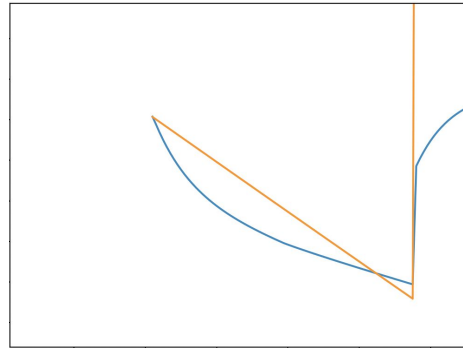
(array([5795.37993376]), (4.312862233015471, 0.008998123942075331, 0.8242315014070988, 47234.95057604644, 0.5715385482453891, 8952.559374074575, 0.12797498520881034, 43311.63512766412))

R\_2

C\_2

R\_3

C\_3



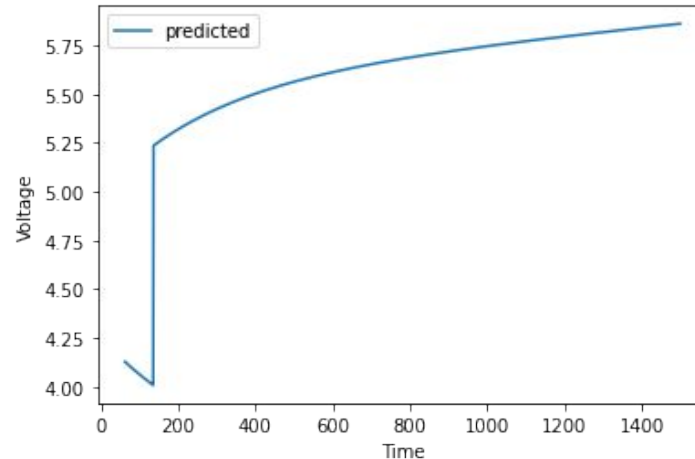
# GENETIC ALGORITHM

## > Generations of solutions

```
====gen23 best solutions====
```

	E_m	R_0	R_1	C_1
(array([795.03570759]),	(4.569385234542004,	0.03637249235303941,	0.6183143238353149,	33092.443575661164,
0.5816582703919699,	28911.1598408088,	0.8903371089946931,	16532.519996790288))	

R_2	C_2	R_3	C_3
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# GENETIC ALGORITHM



- > Able to get approximate solutions using Genetic Algorithm
- > Best solution tricky to obtain but not impossible
- > Genetic Algorithm combined with Neural Networks will give best optimal solution but **very** slow in computation
- > Efficient optimization algorithm with endless opportunities to solve difficult problems

# Sparse Identification of Non-Linear Dynamics (SINDy)



A model discovery technique which identifies non-linear relationships in dynamical systems

Algorithm is as follows:

Input: library of functions of our choice

Output: weights of the functions we gave as input

A typical SINDy model finds the relationship between a state vector (with its combinations) and its derivative

In our model, we tried to equate different combinations of input features to an output vector



# SINDy Algorithm

- Define the current input and true voltage output
- Define the sample time interval
- Input the library of functions
- Call the function used to calculate the parameter vector
- Calculate the predicted voltage value using the parameter vector and library of functions
- Plot the true and predicted voltage vs time
- Calculate the mean squared error

```
l=1439;
b=zeros(l,1)+10; % input current
t=1:1:l; % sample time interval
t=t';
theta = [b b.*(1-exp(-t/80)) b.*(1-exp(-t/800)) exp(-t/500) log(t)]; % library
lambda=0.023; % regularization parameter
n=1;
Xi = sparsifyDynamics(theta,volt,lambda,n); % calculating weights
syms ti I;
Thetasym=[I I*(1-exp(-ti/80)) I*(1-exp(-ti/800)) exp(-ti/500) log(ti)];
Thetasym=Thetasym';
vpredic=transpose(Xi)*Thetasym;
J=0;
vpredic=1:1:l;
vpredic=vpredic';
for i=1:l
    vpredic(i)=subs(vpredic,[ti,I],[i,b(i)]); % calculating predicted voltage
    J= J + ((vpredic(i)-volt(i))^2/l); % calculating mean-squared error
end
```

# Results

Here is the governing equation we obtained from the algorithm,

$$\text{Voltage} = w_0 + w_1 * I * (1 - e^{-t/80}) + w_2 * I * (1 - \exp(-t/800)) + w_3 \exp(-t/500) + w_4 \log(t)$$

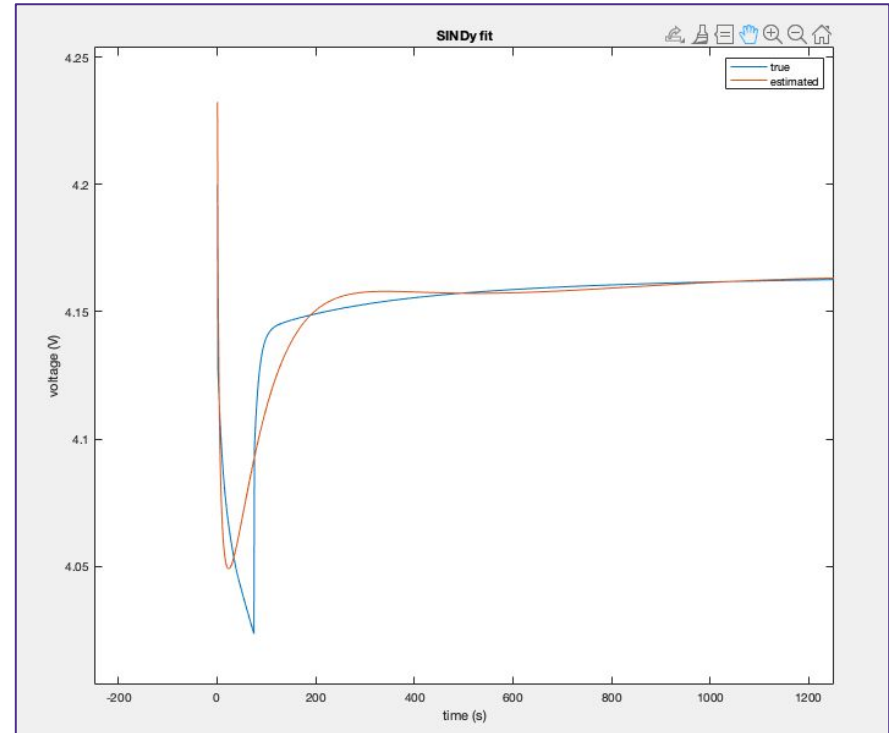
where  $w_0$  ,  $w_1$  ,  $w_2$  ,  $w_3$  ,  $w_4$  are the parameters obtained from SINDy algorithm and  $t$  denotes sample time.

The values obtained are

$$w_0 = 0.41 , w_1 = 0.038 , w_2 = 0.03 , w_3 = 0.08 , w_4 = - 0.089$$

# Model fit and Observations

- The figure on the right shows the fit of the SINDy model
- We can see that the fit is not perfect but follows the pattern and the MSE =  $8.26e-05$  sq volts
- SINDy is not able to capture the entire dynamics probably due to the fact that the input is a constant in this case

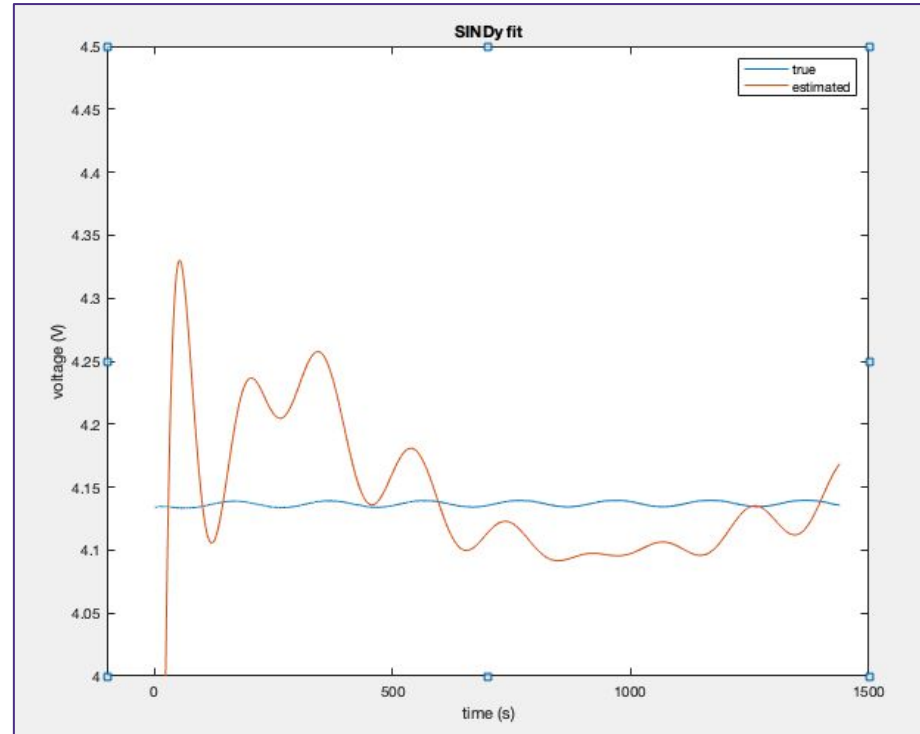


# Validation

- We couldn't obtain the test data for the problem in hand
- So we used MATLAB's System Identification toolbox to generate an 87% fit Hammerstein - Wiener Non-linear model to generate the test data using a sinusoidal current profile of frequency 200 Hz

# SINDy validation

- The figure on the right shows the fit of the SINDy model for test data
- We can see that the fit is not looking great, but tries to follow the pattern and the MSE = 0.0168 sq volts



# Regression with SGD and Levenberg - Marquardt Algorithm

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- Stochastic Gradient Descent is an extension of Gradient Descent which computes gradient of random input data points, thus reducing computational expense
- Levenberg - Marquardt algorithm is a combination of gradient descent with Newton - Raphson method which calculates the hessian matrix in addition to the gradient matrix
- Unlike SINDy which finds hidden dynamics, this algorithm requires a governing equation to be input and finds weights

# Algorithm

This is our governing equation (equivalent circuit model):

$$V_{\text{batt}} = \text{OCV} - I * R_0 - I * R_1 * (1 - e^{-t/(R_1 * C_1)}) - I * R_2 * (1 - e^{-t/(R_2 * C_2)}) - I * R_3 * (1 - e^{-t/(R_3 * C_3)})$$

- Here is the basic equation of Levenberg Marquardt algorithm we used with SGD:

$$x_{n+1} = x_n - [ \nabla^2 f(x_n) + \lambda * I ]^{-1} * \nabla f(x_n)$$

where

- $x_{n+1}$  - vector of parameters at the next time step
- $x_n$  - vector of parameters at the current time step
- $\nabla^2 f(x_n)$  - hessian matrix
- $\lambda$  - Learning rate
- $\nabla f(x_n)$  - gradient

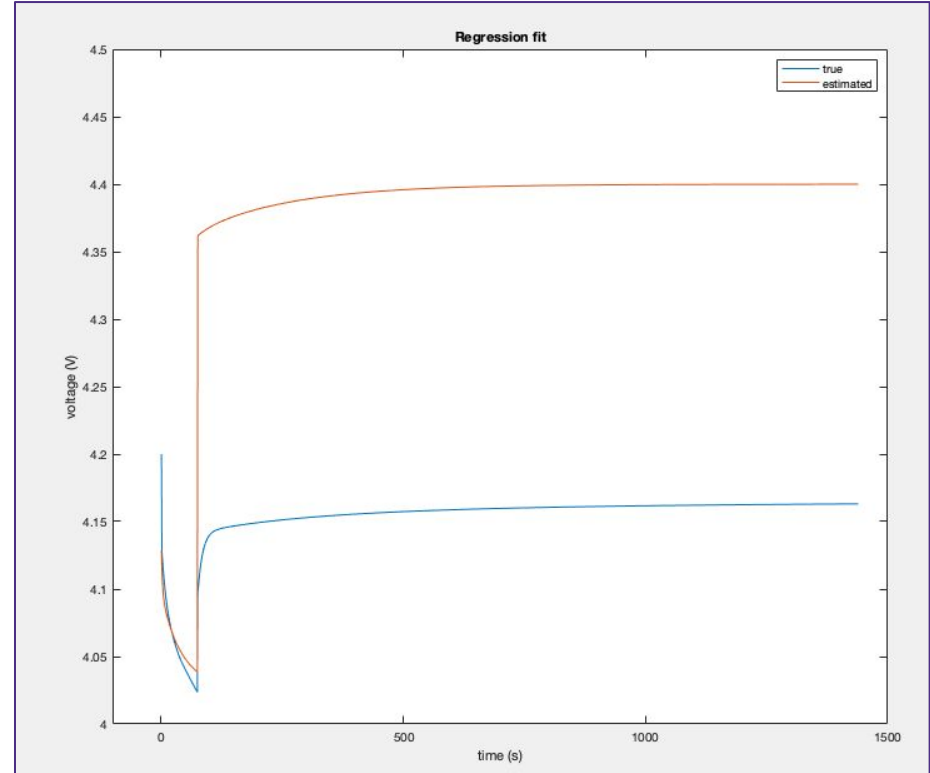
# Algorithm

- Define the governing equation, current input and true voltage output
- Define the cost function
- Define the mini-batch of random points for stochastic gradient descent algorithm
- Calculate the Hessian and gradient
- Execute the Levenberg-Marquardt algorithm
- Use the obtained parameters to calculate the predicted voltage output
- Plot the true and predicted output with time
- calculate the mean squared error



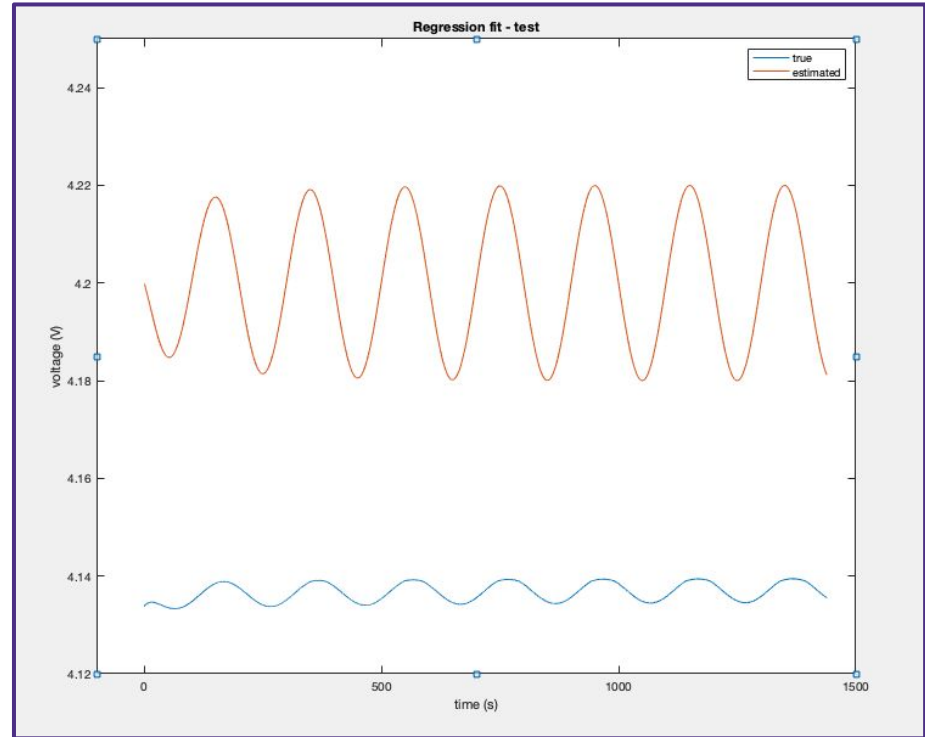
# Results and Inference

- This is the best fit we could get with MSE = 0.0533 sq volts
- Fit largely depends on the initial guess of weights
- Initial guess is done using known physics of the system
- Even a full gradient descent doesn't provide better results because of the local minima issue



# Regression validation

- The figure on the right shows the fit of the Regression model for test data
- The fit follows the same sinusoidal pattern as the true test data with an offset and MSE = 0.0041 sq volts



# Gaussian Process Regression

What is the objective?



- It allows us to make predictions about our data by incorporating prior knowledge.
- Fully Specified by training data, mean and covariance functions
- Covariance is given by “kernel” which measures distance of inputs in kernel space

# Gaussian Process Regression

## Definition

- Given Inputs ( $\mathbf{X}$ ) and output ( $\mathbf{y}$ ):

$$\mathbf{D} = \{(\mathbf{x}_1, y_1), (\mathbf{x}_2, y_2), \dots, (\mathbf{x}_n, y_n)\} = (\mathbf{X}, \mathbf{y})$$

- We presume that there is some structure in our data  $\mathbf{D}$ . GPs model the output as a noisy function of the inputs:

$$y_j = f(\mathbf{x}_j) + \varepsilon; \text{ where } \varepsilon \sim N(0, \sigma_n^2)$$

- Given a set of inputs ( $\mathbf{X}$ ), GP models the outputs ( $\mathbf{y}$ ) as jointly gaussian:

$$P(\mathbf{y} | \mathbf{X}) = N(\underbrace{m(\mathbf{X})}_{\text{Mean}}, \underbrace{K(\mathbf{X}, \mathbf{X})}_{\text{Kernel}} + \underbrace{\sigma_n^2 I}_{\text{Noise}})$$

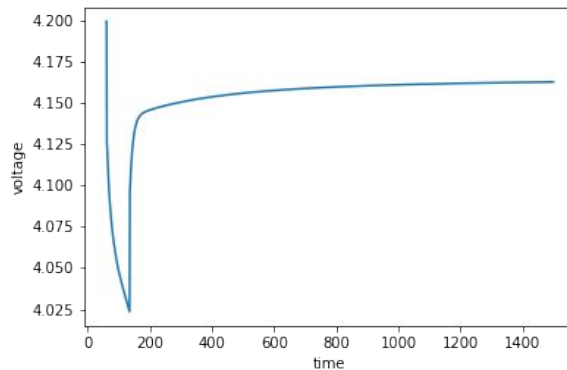
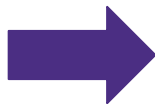
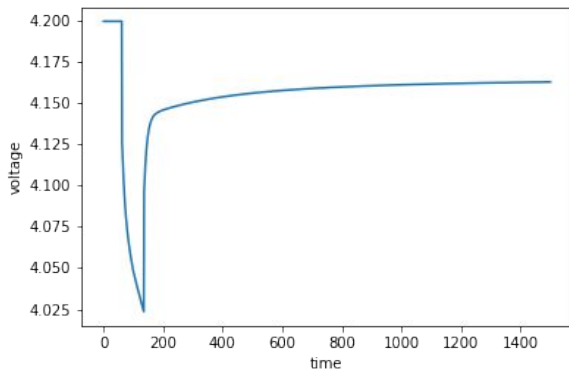
Mean    Kernel    Noise

- Usually we assume prior mean to be zero
- Covariance matrix  $\mathbf{K}$  is defined through kernel function. Example: Squared exponential or RBF, Marten, Periodic exponential

# Gaussian Process Regression

## Training Data

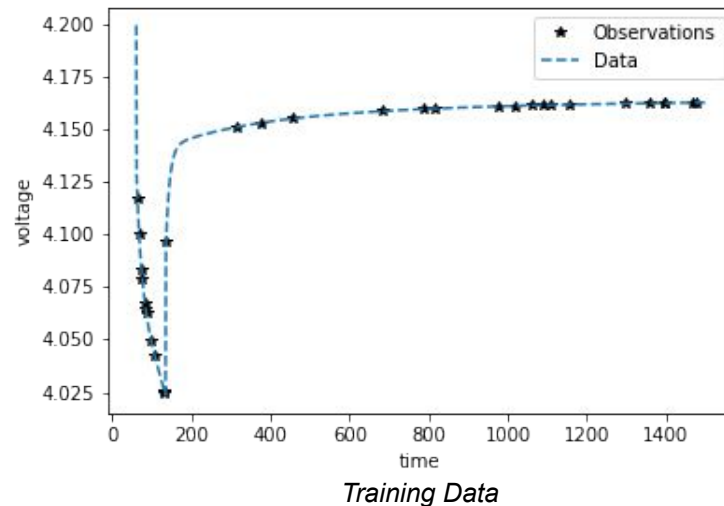
- For our system, current and time are the input parameters. Our Input matrix  $X \in \mathbb{R}^{n \times 2}$  and our output  $V \in \mathbb{R}^n$  where  $n$  is the size of the dataset.
- We are working on the first discharge pulse,  $n = 1500$ . As from  $t = 0s$  to  $t = 61s$  the battery is left ideal, we remove this portion from our analysis.



# Gaussian Process Regression

## Training Data

- First, Sample 30 points from the data randomly.
- Added a bias so that there are some points sampled from the region when the battery is being discharged
- These samples are our observations.
- We take all of the data as our test points
- Setup up input matrix  $X_{\text{train}}$  and  $X_{\text{test}}$  by vertically stacking  $[i(t), t]$



# Gaussian Process Regression

## Setting up Kernel Function

- To generate Covariance matrix, we evaluate kernel pairwise on all the points

$$\mathbf{K} = \begin{pmatrix} k(x_1, x_1) & \dots & k(x_1, x_n) \\ k(x_2, x_1) & & \\ \vdots & k(x_i, x_i) & \vdots \\ k(x_n, x_1) & \dots & k(x_n, x_n) \end{pmatrix}^{**}$$

### Radial Basis Function kernel (RBF kernel)

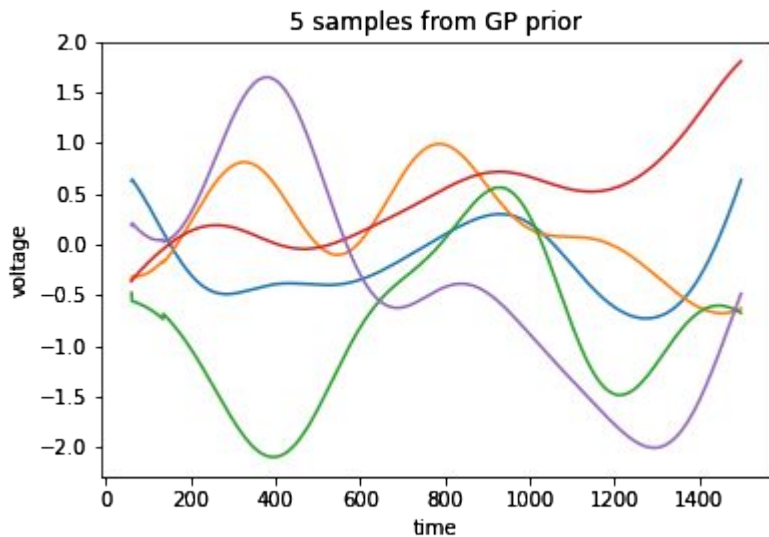
- $\mathbf{K} = \sigma \exp(- ||x - x' ||^2 / 2 \ell^2)$ , where
  - $\ell$  is length scale
  - $\sigma$  is signal variance

### Marten Kernel

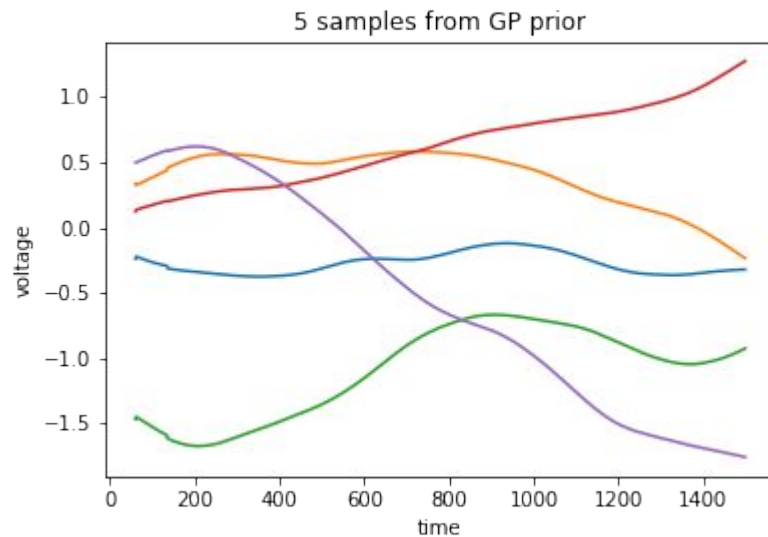
- $\mathbf{K}_{\nu=3/2} = (1 + \sqrt{3} r / \ell) \exp(- \sqrt{3} r / \ell)$ 
  - $r$  is euclidean distance
  - $\ell$  is length scale

# Gaussian Process Regression

## Functions Sampled from Prior



*Prior using RBF kernel*



*Prior using marten kernel*



# Gaussian Process Regression

## Prediction

- Recall, GP models the output  $y$  as jointly gaussian.
- Given training data set  $D = \{(\mathbf{x}_1, y_1), (\mathbf{x}_2, y_2), \dots, (\mathbf{x}_n, y_n)\} = (\mathbf{X}_{\text{train}}, y_{\text{train}})$  and test data set  $(\mathbf{x}_*, y_*)$ ,

$$P(y_{\text{train}}, y_* \mid \mathbf{X}_{\text{train}}, \mathbf{x}_*) = N(\mu, \Sigma)$$

- Condition on  $y$ : As the outputs are jointly gaussian, marginalisation on the known outputs would also result in a gaussian.

$$P(y_* \mid \mathbf{x}_*, y_{\text{train}}, \mathbf{X}_{\text{train}}) = N(\mu_*, \Sigma_*),$$

where  $\mu_*, \Sigma_*$  are the mean and covariance of the functions and are given by:-

$$\mu_* = K_*^T (K + \sigma_n^2 I)^{-1} y_{\text{train}}$$

*mean of predicted functions*

$$\Sigma_* = K_{**} - K_*^T (K + \sigma_n^2 I)^{-1} K_*$$

*Covariance of predicted functions*

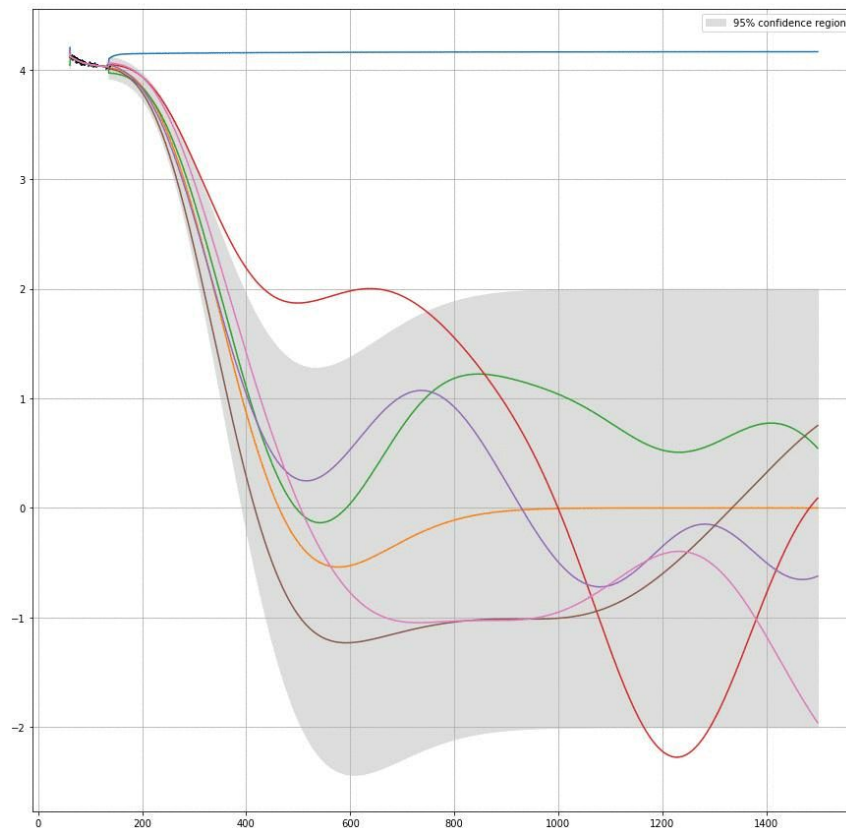
$K_*$  = captures relation between training points and test points

$K_{**}$  = captures relation among test points

# Gaussian Process Regression

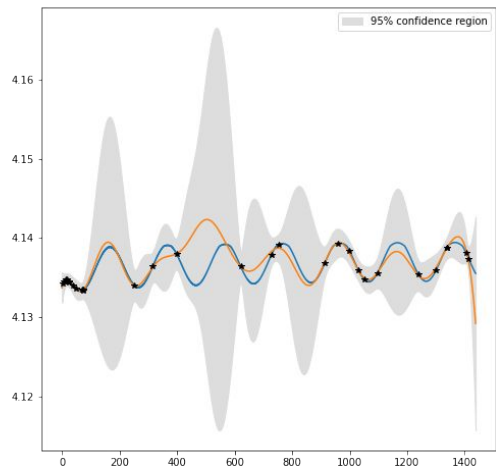
## Prediction

- Prediction using RBF kernel

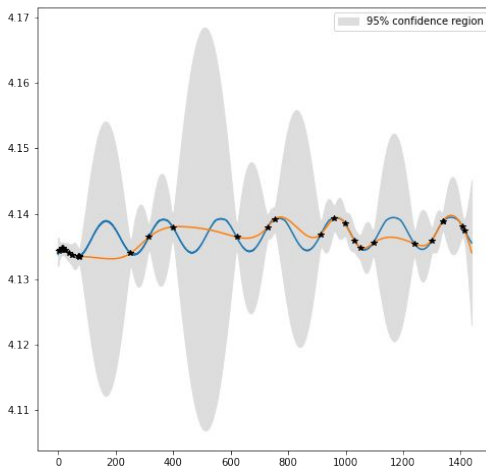


# Gaussian Process Regression

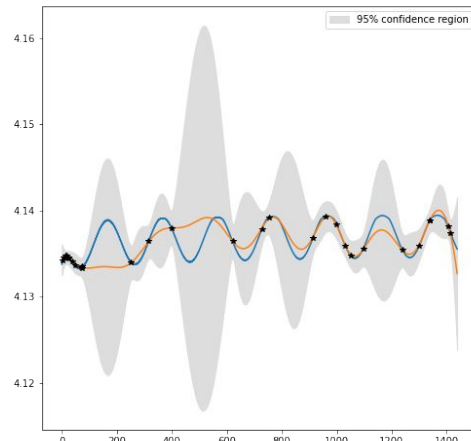
## Validation



*RBF kernel*



*Marten kernel  $\nu = 1.5$*



*Marten kernel  $\nu = 2.5$*

# Conclusion



- In this project we explored 4 different data driven techniques for system identification or model discovery
- **SINDy** gives information about the system dynamics and is helpful to find the nonlinear dynamics
- **Genetic Algorithm** can be used to solve search and optimization problems efficiently
- **Regression Stochastic Gradient** is a powerful method for gray box model discovery
- **Gaussian Process Regression** gives a distribution over a set of functions. They allow us to make predictions about our data using prior knowledge.

# References



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