# Chapter 16 State of Energy Estimation of Li-Ion Batteries Using Deep Neural Network and Support Vector Regression



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**Abstract** Efficient management of the power and energy output of a high voltage battery pack requires a precise estimation of the State of Energy (SOE). For the accurate estimation of SOE, this work presents two data-driven methods as Deep Neural Network (DNN) and a regression model, i.e. Support Vector Regression (SVR). The effectiveness of the SOE estimation was compared, analysed, and studied through these models under similar conditions. For performance enhancement of estimation, a modified algorithm based on the grid search of optimized hyperparameters was proposed and evaluated in both the models. For training of the model at subsequent thermal ranges, two case studies were performed using US06, UDDS, LA92, and HWFET drive cycles and at four different temperature levels (-10, 0, 10, and 25 °C), for each cycle. The results indicate that the DNN method has provided enhanced performance for State of Energy Estimation as compared to the regression models of ML, i.e. SVR. This work highlights the prevailing challenges in the industry and proposes the potential recommendation for Battery Management System (BMS) development and SOE estimation in next-generation EV applications.

**Keywords** The State of Energy · Deep neural network · Support vector regression · Hyperparameters optimization · Regression

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<sup>©</sup> The Author(s), under exclusive license to Springer Nature Singapore Pte Ltd. 2022 299 I. Ali et al. (eds.), *Computational Modelling in Industry 4.0*, https://doi.org/10.1007/978-981-16-7723-6\_16

#### 1 Introduction

With the growing problems such as the depletion of energy resources and the problem of global warming caused by use of internal conventional engines-based vehicles, electrical vehicles (EVs) have attracted the very high attention of people (Tie and Tan 2013; Xia et al. 2017). In recent years, a lot of development has taken place in the electrochemical energy storage system. Different energy storage systems have been proposed for the use in EVs such as Nickel/Metal Hydride (NiMH) battery, Lithium ion (Li-ion) battery, Fuel cells, ultra-capacitors, etc. (Iclodean et al. 2017). With the advancement and improvement in the properties of Li-ion batteries such as energy density, low self-discharge rate, long cycle life, and safety performance, they become widely applicable in EVs, electronics, mobile devices, etc. (Lu et al. 2013). A Li-ion battery module consists of an array of Li-ion cells. Temperature management is one of the issues of the Li-ion batteries which affect performance and safety (Gandoman et al. 2019; Kumar et al. 2020). The literature shows several solutions to improve the cooling effects by optimized battery layout (Qian et al. 2019), Heat Pipe (Lu et al. 2020), Phase Change Materials (Jaguemont et al. 2018), modular and mixed solutions (Cicconi et al. 2020). This paper aims to estimate the level of State of Energy (SOE) considering the achieved cell temperature with the operation parameters, i.e. voltage and current. The estimation of the cell state parameters such as State of Charge (SOC), Open Circuit Voltage (OCV), State of Energy (SOE), etc., is very important for the accurate functioning of the system and for the long life of batteries. These state parameters are also important in the estimation of driving range of electric vehicles which is also one of the complex issues as studied by Ronan German et al. (2020). However, the correct estimation of these state parameters is quite complex and difficult due to the non-linear behaviour of the electrochemical processes in Li-ion cells (Hafsaoui and Sellier 2010).

Traditionally, SOC estimation is performed for the identification of residual energy and the protection of batteries from being overcharged and discharged. From the recent literature review, it is well shown that different researchers had used a wide variety of techniques for the accurate estimation of SOC such as current integral method (Ng et al. 2009), proportional-integral (PI) method (Xu et al. 2014), electrical model-based method (Plett 2004b, c; He et al. 2013; Zhong et al. 2014), Sliding-mode observer based (Kim 2006, 2010), Kalman filter-based algorithms (Plett 2004a, b, c; Xu et al. 2012; Xiong et al. 2014), and neural network model (Kang et al. 2014) methods. With the increasing demand of Li-ion batteries for different applications, there is increased demand of use of Battery Management System (BMS) as well. Therefore, it is more and more important to accurately analyse the battery states for the proper functioning of batteries. But Liu et al. (2014) reviewed several disadvantages in the use of SOC for the estimation of residual energy. Because SOC is defined as the ratio of available capacity to the maximum stored charge in the battery, i.e. nominal capacity, from which it can be seen that there is no representation of the State of Energy. SOC clearly gives the information only about the residual

capacity, not about the energy. That's why there is the need to estimate SOE independently. SOE is defined as the available energy to the maximum stored energy in Li-ion battery. Some researchers (Shen 2007; Hausmann and Depcik 2013; Waag and Sauer 2013; Zheng et al. 2013) used residual available capacity instead of SOC for the estimation of SOE. Waag and Sauer (2013) used the estimation of battery electromotive force for the identification of the battery capacity and SOC. While Shen et al. (2007) defined state of available capacity of battery, instead of SOC for the estimation of battery residual capacity. These are some of the few works which used the capacity instead of SOC in the study. It is also important to study the effect of discharge current and temperature, since at the same SOC; SOE can be different because discharge efficiency is dependent on discharge current and temperature (Liu et al. 2014). Few researchers worked in this direction to understand the working of Li-ion batteries, as Wang et al. (2014a, b) show the study on the estimation of electronic conductivity of LiFePO4 cells at different temperatures to understand the low-temperature electrochemical performance at carbon-coated and uncoated cathodes. Yi et al. (2013) developed a model for the study of behaviour of Li-ion battery of temperature dependence on discharge in low ambient temperature. These studies prove that for the accurate estimation of SOE, it is necessary to consider the effect of discharge current and temperature.

In recent years, different researchers developed different techniques for the estimation of SOC and also proved well with the results. But SOC is different from SOE because SOE is the product of the residual battery capacity per OCV. The trend of variation of SOE is different from SOC. Similar to the estimation of SOC, few researchers developed systematic methods for the SOE measurement (Mamadou et al. 2012, 2019; Liu et al. 2014). In these studies, direct evaluation techniques were shown for the residual energy of battery with the consideration of battery discharge states. Some other researchers such as Stockar et al. (2011) and Kermani et al. (2011) presented the method of power integral for the estimation of SOE. But these studies were not found very accurate for the SOE measurement due to the measurement noises in the current and voltages of battery. Wang et al. (2014a, b) proposed the joint estimation technique of SOC and SOE to minimize the negative effect of power integral method. But in this method SOE estimation accuracy depends on the SOC estimation accuracy. To overcome this issue, Zhang et al. (2015) proposed the method of model-based joint estimation of SOC and SOE.

From the literature reviews, it can be seen that there is still a lot of possibility of research for the estimation of SOE. Since in this area, research has not been done as far as compared to the SOC estimation. Also, it was found out that SOE estimation is somewhat hard to estimate using mathematical relations and equations. So, there is a need to develop a method which is less complex and computationally efficient to be employed in real BMS condition. Therefore, this paper presents the two datadriven approaches for the estimation of SOE such as a regression model, i.e. SVR and DNN. A modified DNN-based method is shown for the accurate estimation of SOE. The modified method consists of the two simultaneous processes, i.e. first the optimization of hyperparameters and then the DNN model development which will be used for the prediction. In the neural networks, the critical task is the optimization of hyperparameters which is accurately done using a grid-based approach. In the proposed grid-based approach, grids were developed based on the Root Mean Square Error (RMSE), training time, and hyperparameters such as the number of neurons and number of layers. Optimized values of the number of neurons and number of hidden layers were chosen when minimum RMSE with suitable training time was recorded. The same process was used in the development of regression model using SVR for the estimation of SOE. Grid-based approach was used in the modified algorithm of SVR model for the searching of optimized hyperparameters such as regularization parameter (C) and gamma ( $\gamma$ ).

The main aim of this study is to develop an efficient model for the estimation of SOE and to overcome the shortcomings of regression model. For the implementation of such task, in this study two data-driven methods were used for the development of SOE estimation model, DNN and SVR. The results obtained from the DNN model supports that it is more efficient in the estimation as compared to the SVR. The results shown in Sect. 4 clearly state that the DNN model is fully efficient in overcoming the shortcoming of regression model. DNN model shows significant results in the prediction of SOE at different drive cycles and at different temperatures as well.

#### 2 Proposed Methodology for SOE Estimation

The SOE of battery can be defined by Eq. (1) as the ratio of residual energy to the maximum available energy, where residual energy is denoted by  $E_{res,k}$ , maximum energy by  $E_M$ , battery energy efficiency as  $\eta_e$ , battery terminal voltage and current at kth time instant as  $V_{t,k}$  and  $I_k$ ,  $T_s$  is the time sampling period. The methodological approach for the estimation of SOE is described in Fig. 1. The approach is focused on the accurate estimation of SOE followed by the selection of optimized hyperparameters for the respective estimation learning models. As shown in Fig. 1, raw data was collected from the publicly available dataset (Kollmeyer et al. 2020). The data for SOE was processed from this dataset using the Eq. (1). Data cleansing was performed in the next step such as resampling of the dataset to 1 Hz frequency and the removal of errors in data (such as Voltage, Current spikes). After the data cleansing, the further dataset was developed according to the respective case studies of this research. In this study, two estimation learning model was selected, i.e. DNN and SVR. Optimization of hyperparameters was the main step after the choosing of the training dataset, validation dataset, and the selection of model. Hyperparameters are responsible for estimation accuracy, computation time, and computational speed. The processes of estimation in both models are shown in detail in Sect. 2.1 and 2.2. The performance evaluation was done on the basis of RMSE and MAE. The evaluation accuracy is discussed in the result section, i.e. Sect. 4.

$$SOE_{k+1} = \begin{cases} SOE_k - \frac{E_{res,k}}{E_M} \\ SOE_k - \frac{\eta_e V_{i,k} R_J_s}{E_M} \end{cases}$$
(1)



Fig. 1 The proposed methodological approach for SOE estimation

#### 2.1 DNN Architecture

DNN is part of the family of machine learning methods based on artificial neural networks (ANN). In the present time, DNN has wide applications, there are many examples available where DNN architecture shows good response over the conventional algorithms such as DNN showed the level of accuracy more the human level such as in 2015 Microsoft Research's deep neural algorithm won the ImageNet challenge with the error of about 3.57% (He et al. 2016).

DNN is considered fast and efficient in comparison to the conventional methods for the estimation of SOE, since in conventional method expertise is required in battery chemistry and to mathematical model the battery behaviours. On the other hand, DNN doesn't involve such complexity for the SOE estimation. DNN consists of three layers, i.e. input layer, hidden layer, and output layer. The input and output layer remains fixed while the hidden layer can be varied to find out the optimum value where minimum error can obtain. Along with the layers, nodes or it can be said as neurons are also present in the respective layers to capture the non-linearity between the system input and output. The optimum value of nodes is also required to be found with the hidden layers for the development of the efficient and computational fast model.

DNN learns the battery functioning by mapping the input parameters with the output parameters. In this study for the creation and validation of DNN model a

publicly available dataset was used (Kollmeyer et al. 2020). The dataset consists of battery parameters of LG 18650HG2 Li-ion battery at four different drive cycles such as US06, UDDS, HWFET, LA92, and all the data of drive cycles are available at four different temperature set, i.e. 10, 0, 10, and 25 °C. Battery parameters available in the dataset are current, voltage, temperature, and SOE with time step of 1 s. For the development of the DNN model, vector of battery parameters was provided to the input layer where parameters were current, voltage, temperature, average current, and average voltage, while the output layer was having the single parameter, i.e. SOE. The model was created in such a way that it can effectively map the variation in input parameters with the variation in output, i.e. SOE and by using this learning method, it will be able to predict the SOE for any new input parameters value. Mathematically the input vector can be represented as— $X(t) = [I(t), V(t), T(t), I_{avg}(t), V_{avg}(t)]$  and the output can be represented as—Y(t) = SOE(t), where I(t), V(t), T(t),  $I_{avo}(t)$ ,  $V_{avo}(t)$ , and SOE(t) denote the current, voltage, temperature, average current, average voltage, and estimated SOE at time step t. The input vector X(t) fed to the input layer to map with the SOE(t) values at the output layer. Figure 2 illustrates the architecture of deep neural network with input, output, and hidden layers along with the neurons in each layer. It is showing the process of mapping of observables with the desired output. The computational speed of offline training of DNN is fast because to capitulate the output Y(t), input vector X(t) perform certain matrix multiplication whereas other strategies are not so computational fast since it involves partial differential equations. The hidden layer activation is represented by Eq. (2). Since DNN is matrix-based,



Fig. 2 Architecture of 2-layer DNN with representation of input, hidden and output layer

there are certain variables which need to be defined such as  $w_{j,k}^{L}$  (which denotes the weights connection between neuron j in layer L-1 and neuron k in layer l),  $b_{k}^{L}$  (which denotes the bias), and  $h_{k}^{L}$  (which denotes the activation, respectively, of neuron k in layer L).

$$h_k^L = \sigma\left(\sum_k \left(w_{j,k}^L h_k^{L-1}(t) + b_k^L\right)\right)$$
(2)

Reading Eq. (2),  $\sigma$  is the activation function. Due to simplicity in training and testing, the non-linearity or activation function used for hidden layer is Rectified Linear Unit (ReLU) which can be given by Eq. (3).

$$\sigma(\mathbf{x}) = \begin{cases} 0 for \mathbf{x} < 0\\ x for \mathbf{x} \ge 0 \end{cases}$$
(3)

Similarly, at the output layer, SOE can be calculated using Eq. (4).

$$SOE(t) = \eta \left( \sum_{k} \left( w_{j,k}^{L} h_{k}^{L-1}(t) + b_{k}^{L} \right) \right)$$
 (4)

where L is the last hidden layer of network and  $\eta$  is the activation function for output layer which was chosen as Sigmoid function and can be represented by Eq. (5).

$$\eta(x) = \frac{1}{(1 + \exp(-x))}$$
 (5)

#### 2.1.1 DNN Hyperparameters and Its Optimization

Effective modelling of DNN requires selection of optimum values of hyperparameters because the computational speed and performance of model depends upon these hyperparameters. There are number of hyperparameters available such as hidden layers, neurons, learning rate, optimization algorithm, activation functions, etc. But the optimization of all the hyperparameters at the same time will not be a timeeffective approach; therefore the most important hyperparameters were optimized in this study which was number of hidden layers and number of neurons in hidden layer. Rest of the hyperparameters were selected based on the literature survey such as learning rate was kept 0.001, optimization algorithm was chosen as Adam due to its good performance, ReLU activation function was used for hidden layers and Sigmoid function for output layer. Optimization algorithm developed in this study for the hidden layers and neurons selection is shown in Sect. 2.1.2. Grid-based approach was used for developing the optimization algorithm and US06 drive cycle at 25 °C was used for the validation in this optimization technique. Range of values was chosen both for hidden layers and neurons, to search the combination where root mean square error (RMSE) minima will occur. The range of values checked for hidden layers was from 2 to 8, while the neurons set of values were (2, 4, 8, 16, and 32). Table 1 shows the values of RMSE and training time for the different combinations of hidden layer and neurons. From Table 1, it can be seen that the minimum RMSE occur at the combination of 3 hidden layers and 16 neurons with training time as 775.7467 s. Therefore, keeping in mind about the computational efficient concept, combination of 3 hidden layers and 16 neurons in each hidden layer was chosen for the further study. Table 2 shows the list of all the optimized hyperparameters used in this study.

Layers	2	3	4		5	6	7	8
Neurons								
Root mean Square Error (RMSE)								
2	0.173742	0.161347	0.286389		0.143908	0.145942	0.286164	0.287129
4	0.098428	0.110402	0.285939		0.027262	0.08073	0.170202	0.051682
8	0.089132	0.038056	0.030829		0.022372	0.017867	0.016731	0.015806
16	0.083765	0.014997	0.019165		0.017265	0.018367	0.016447	0.015747
32	0.040811	0.036811	0.031811		0.028821	0.021734	0.0199	0.017876
Training 2	Time (secon	ds)						
2	640.9791	694.9241	719.4053	841.	239	895.1712	939.1342	951.2096
4	568.7269	751.9329	582.862	770.	4256	831.8109	864.1268	831.5526
8	582.5267	774.7212	592.3911	769.	8025	893.7469	890.4075	913.9425
16	586.9675	775.7467	583.1441	657.	4654	815.7815	871.2201	855.9458
32	590.2175	800.4176	608.9079	686.	7024	976.5569	1007.908	852.5247

 Table 1
 Grids showing the values of RMSE and training time for different combinations of hidden layers and neurons

Table 2Optimized valuesselected for thehyperparameters of DNN

Hyperparameters	Optimized values
Number of hidden layers	3
Number of neurons	16
Optimization algorithm	Adam
Activation function for hidden layers	ReLU
Activation function for output layer	Linear
Learning rate	0.001

#### 2.1.2 DNN Training Strategy

This section gives the information regarding the strategy used for the training of DNN, with the insight about incorporation of optimization technique used for hidden layer and neurons along with the training. In this study, before starting the training of the model optimized value for number of hidden layers and neurons were found and the selection procedure is shown in Sect. 2.1.1. Appendix A.1 shows the modified algorithm used in this study for the development of effective DNN model. Algorithm starts with the importing of dataset used for the training. Dataset used in this study were at the different drive cycles such as US06, UDDS, HWFET and LA92. These drive cycles have the values for different parameters such as Current, Voltage, Temperature, Average Current, Average Voltage, and SOE at four different temperature set, i.e. -10, 0, 10, and 25 °C. Input parameters that will be fed into the input layers were put in a vector X and output parameter, i.e. SOE in Y. Since neural network have 3 layers, input, output, and hidden layers, input and output layers were ready as X and Y but number of hidden layers and number of neurons in each layers were still needed to be found and optimized. Therefore for finding the optimized values, range of values for hidden layer and neurons too were decided. RMSE was calculated for the different combinations of these values using the grid-based approach. Table 1 shows the RMSE and training time values for the different combinations. Finally, after the selection of optimized values of number of hidden layers and neurons the modelling of DNN begins. Optimization of these hyperparameters are utmost important because performance and efficiency of model depends on these parameters. In this study, the activation function of hidden layer was chosen to be ReLU, since from the literature study it was found to be most efficient among the other alternatives. The model used in this study updates the weights with the Adam optimization technique with learning rate 0.001. Mean Squared Error (MSE) loss function and Mean Absolute Error (MAE) metrics was chosen during the training of model. Performance analysis was done using the RMSE, as it can evaluate the percentage deviation of predicted values from original values. Training was done using the computer system with specifications of 4th generation core i5 processor and intel HD graphics. In all the cases of this study, model was trained for 1000 epochs to minimize error and to improve the model efficiency. Models were developed for 2 different cases, as in Case 1 model was trained with 3 drive cycles, i.e. UDDS, HWFET, and LA92 at all temperatures set and then performance evaluation was done using US06 drive cycle at different temperatures. In Case 2, the accuracy of model was analysed in mapping the temperature effect. All drive cycles were used for training but considering only three temperature set, i.e. -10, 0, and 25 °C. In this case evaluation was done using three drive cycles as UDDS, LA92, and HWFET at 10°C. The main purpose of considering these case studies were to evaluate the performance of model at some new temperature and new drive cycle which was not considered during the training. Results and comparison of these 2 cases are shown in Sect. 3. The performance of all the models was evaluated on the basis of MAE and RMSE error metrics.

#### 2.2 SVR Architecture

In machine learning, Support Vector Machines (SVM) is the supervised learning models which analyse data for the classification and regression analysis. This study focuses on the SVR analysis technique. The functioning of SVR is in the way that it makes a complex relationship between input parameters and single output by optimizing the objective function, whose values get minimized when the predicted value from model approaches the target or original output value. For this case also the input parameters remain same as it was used for DNN, i.e. current, voltage, temperature, average current and average voltage. The major task of SVR is to correctly and precisely map the input parameters with the output value, i.e. SOE. For this study, it is not possible to have a linear relationship between the input and target values. Therefore, for this type of circumstance input values get linearly mapped with the output values in higher dimensional plane and the functions used for these transformations are known as Kernels (Müller et al. 2001). The available kernels functions in the SVR technique are sigmoid function, polynomial function, Radial Basis function (RBF), etc. Among the other available kernel function and based on the nature of the relationship between input and output parameters, RBF kernel was the most suitable for this study (Schölkopf et al. 1997). The architecture of the SVR is shown in Fig. 3, where an insight about the mapping of input and output parameters can be understood easily. From Fig. 3, it is shown that  $K(x, x_n)$  is the output of the nth hidden node for the input X, also it is the mapping of input vector x and support vectors by precisely choosing kernel function (Chen and Yu 2007).

In the SVM technique, the two regression methods are available, i.e.  $\epsilon$ —SVR and  $\nu$ —SVR. Both methods differ from each other in a way that in  $\epsilon$  – SVR methods



Fig. 3 U—SVR architecture

there is no control on the number of data vectors from input dataset that become support vectors. While in  $\upsilon$ —SVR method, there is the control on the limit of error tolerance where  $\upsilon$  represents the upper bound on the fraction of error from training dataset and the lower bound on the fraction of support vectors. This control on limit of error tolerance can regulate by the use of regularization parameter C. For this study, authors find  $\upsilon$ -SVR method more suitable since error tolerance in the prediction can be controlled.

#### 2.2.1 SVR Hyperparameters and its Optimization

Similar to DNN modelling, the SVR modelling also involves hyperparameters handling and its optimization. In the SVR analysis the main hyperparameters which affect the performance and efficiency of model are gamma ( $\gamma$ ), C, kernel function, and v. The y parameter shows the range of influence of input training data samples. Simply, it can be defined as the inverse of radius of influence of support vectors selected by model from the dataset. C is the regularization parameter which controls the trade-off between achieving the low training and low testing error. Kernel function is responsible for making the required relationship between input and output parameters. Due to the nonlinear relationship between the input and target parameters of this study, RBF was chosen as the kernel function. The use and functioning of  $\upsilon$  is already explained in Sect. 2.2, it should be in the interval of (0, 1]. As the optimization of all the hyperparameters at the same time cannot be the efficient way to do the required study, therefore the default value of  $\upsilon$  was chosen as 0.5, RBF kept as the kernel function, and then the optimum value was selected for the  $\gamma$  and C using the optimization technique similar to the DNN case. The grid-based approach similar to the one shown in Sect. 2.1.1 for the optimization of DNN hyperparameters was used for the optimization of C and  $\gamma$ . The range of values decided for the regularization parameter, C was -0.01, 0.1, 1, 10 and 100. While three values were chosen for gamma parameter for iterating with the combination in C, i.e. 0.1, 1, and 10. Table 3 shows the values of RMSE and training time for the different combinations of C and  $\gamma$ , validation drive cycle, i.e. US06 at 25°C was used for the testing. From these grids one combination of C and  $\gamma$  was chosen at which RMSE was found to be minimum and the training time was very optimum. The combination which satisfied these criteria was C as 0.1 and  $\gamma$  as 10. The RMSE obtained at this combination was 0.016497, which was minimum among rest of the grid combination and training time was 698.3137 s. All the parameters which were used in this modelling and their respective selected optimum values are shown in Table 4.

#### 2.2.2 SVR Training Strategy

One of the most important parts of modelling in machine learning is the training strategy. This section will give the detail insight to the reader about the modified  $\upsilon$  – SVR algorithm with the simultaneous optimization of hyperparameters. The

С	0.01	0.1	1	10	100			
Gamma (y)								
Root Mean Squar	Root Mean Squared Error (RMSE)							
0.1	0.099553	0.07286	0.065908	0.049061	0.040832			
1	0.049851	0.03734	0.024061	0.017731	0.018968			
10	0.031577	0.016497	0.017534	0.017223	0.007121			
Training Time (seconds)								
0.1	667.7578	655.6799	671.5883	432.3391	1165.202			
1	631.336	651.5541	407.5546	1106.363	7541.776			
10	616.2116	698.3137	1557.31	11363.25	39771.38			

Table 3 Grids showing the values of RMSE and training time for different combinations of C and  $\gamma$ 

Table 4         Optimized values           selected for the         hyperparameters of SVR	Hyperparameters	Optimized values	
	С	0.1	
	γ	10	
	υ	0.5	
	Kernel function	RBF	

modified algorithm used for this method is shown in Appendix A.2. The details of the optimization of C and  $\gamma$  are already shown in Sect. 2.2.1. The datasets used for the training in this case are also similar with the DNN training. In this technique also 2 cases were studied such as in first case, three drive cycles, i.e. UDDS, LA92, and HWFET at temperature set of -10, 0, 10, and 25 °C was used for training. In this case the performance was evaluated by using the US06 drive cycle at all the same temperatures from -10 to 25 °C. In case study 2 as that of DNN training explained in Sect. 2.1.2, such that all the four cycles were fed for the training at three temperatures, i.e. -10, 0, and 25 °C. The validation was done using the three drive cycles as UDDS, LA92, and HWFET at 10°C. All these trainings were done using the same kernel function as RBF and nu (v) as 0.5. Since in the SVR technique there is no involvement of epochs, therefore its computation time fully depends on the selection of these hyperparameters. Training was done using the same computer system with specifications of 4th generation core i5 processor and intel HD graphics. The performance of all the cases was studied on the basis of MAE and RMSE error metrics. The detailed comparison among the results of these cases is shown in Sect. 3.

parameters	Cell parameter	Specification	
	Chemistry	Li[NiMnCo]O <sub>2</sub> (H-NMC) / Graphite + SiO	
	Nominal Voltage	3.6 V	
	Charge	1.5A,4.2,50 mA End-Current (CC-CV) Normal	
		4A, 4.2 V,100 mA End-Current (CC-CV) Fast	
	Discharge	2 V End Voltage, 20A Max Continuous Current	
	Nominal capacity	3.0 Ah	
	Energy density	240 Wh/Kg	

#### **3** Processing of Data for Training and Validation

#### 3.1 Data Collection

In this study for the development of data-driven models and for the evaluation, the open source data was used available from Macmaster University (Kollmeyer et al. 2020). LG 18650HG2 battery was used in the testing for the gathering of this data as mentioned in the technical instruction of the open source data. The detailed information of battery used for the testing is shown in Table 5. The data are available for four different drive cycles, i.e. US06, UDDS, LA92, and HWFET tested at four different ambient temperature conditions as -10, 0, 10, and 25 °C. For this study all the required data for these drive cycles were gathered from this source (Kollmeyer et al. 2020) and then further data were prepared according to the cases of this particular study. The training and validation dataset prepared according to the cases as shown in Sect. 3.2.

#### 3.2 Training and Validation Dataset

In this study, four drive cycles were used for the training in different combination as shown in Table 6. Simulation of battery cells by the use of these driving cycles is very time efficient as well as cost efficient. Use of these drive cycles in battery simulation makes it easier to monitor the behaviour of battery cells during real-world driving patterns without doing any extra experimentation on real physical driving vehicle. Driving cycles used in this study were Supplemental Federal Test Procedure

Cases	Testing dataset	Validation dataset
Case 1	UDDS, LA92, and HWFET drive cycle at -10, 0, 10, and 25 °C	US06 drive cycle at -10, 0, 10, and 25 °C
Case 2	US06, UDDS, LA92, and HWFET drive cycle at -10, 0, and 25°C	UDDS, LA92, and HWFET drive cycle at 10 °C

Table 6 Training and Validation dataset used in the study

(US06), Urban Dynamometer Driving Schedule (UDDS), Unified Driving Schedule (LA92), and the Highway Fuel Economy Driving Schedule (HWFET). These cycles are described in Appendix B. As an example, Fig. 4 shows the behaviour of the battery parameters, i.e. current, voltage, and temperature of US06 drive cycle at all ranges of temperature from -10 to 25 °C. Training and validation dataset used in



Fig. 4 The report of the US06 drive cycle in terms of: **a** Current, **b** Voltage, **c** Cell Temperature. This test was done at four conditions of room temperature: (i) -10 °C, (ii) 0 °C, (iii) 10 °C, (iv) 25 °C

different cases for the study in this research are shown in Table 6. Two cases were studied using DNN and then results of DNN were compared by another Machine Learning technique, i.e. SVR. These two cases are already explained in Sect. 2.1.2 and 2.2.2 according to the respective DNN and SVR training.

All the input vectors in each case were normalized before feeding for the training. Normalization technique was used to improve the convergence rate and to remove the negative influence. In this study min–max normalization was used since it retains the original data distribution pattern, only scaled the dataset in the range of [0, 1] as shown in Eq. (6) (Jain et al. 2005).

$$z_{i}^{k} = \frac{x_{i}^{k} - \min(x)}{\max(x) - \min(x)} i \in \{1, 2, \dots, n\}$$
(6)

where x is the input vectors such as I(t), V(t), T(t),  $I_{avg}(t)$ ,  $V_{avg}(t)$ , and n represents the total number of samples of data in the different drive cycles.

#### 3.3 Evaluation Metrics

For the study of performance of trained models, some evaluation metrics were used such as MAE and RMSE. These evaluation or error metrics made the way of the comparison of different machine learning methods and different case study done in these methods in this research paper. These evaluation or error metrics can be represented by the Eqs. (7), (8), and (9).

$$MAE = \frac{1}{N} \sum_{k=1}^{N} (|SOE_k - SOE_k^*|)$$
(7)

$$MSE = \frac{1}{N} \sum_{k=1}^{N} (|SOE_{k} - SOE_{k}^{*}|)^{2}$$
(8)

$$RMSE = \sqrt{\frac{1}{N} \sum_{k=1}^{N} (|SOE_k - SOE_k^*|)^2}$$
(9)

where  $SOE_k$  is the predicted value by model,  $SOE_k^*$  is the actual value obtained from battery testing at time step k and N is the total number of training samples.

### 4 Results and Discussion

As mentioned in earlier section, the input vector fed in to the DNN and SVR can be represented as  $X = [I(t), V(t), T(t), I_{avg}(t), V_{avg}(t)]$  and the output can be represented as Y = SOE(t), where  $I(t), V(t), T(t), I_{avg}(t), V_{avg}(t)$  and SOE(t) denotes the current, voltage, average current, average voltage and estimated SOE at time step t. The drive cycles used for the training and validation are mentioned in Sect. 3.2 and were recorded at sampling frequency of 1 Hz. The following subSects. 4.1 and 4.2 will show the information about the results obtained after the training and validation according to different cases used in this study. Comparative analysis is shown in these sections for the DNN and SVR modelling for the prediction of SOE.

#### 4.1 Case Study 1

In this case, as described earlier that three drive cycles namely UDDS, LA92, and HWFET were fed for training at all temperature sets as -10, 0, 10, and 25 °C. For the evaluation of the trained model US06 drive was used at all the temperature ranges. Comparative study is shown here between DNN and SVR technique of machine learning. Estimation performance of DNN and SVR was compared on the basis of error metrics, i.e. RMSE and MAE. Figure 5 depicts the prediction of SOE using both data-driven methods, i.e. DNN and SVR. It shows the comparative representation of SOE between actual SOE, predicted SOE using DNN, and predicted SOE using SVR. Figure 5a–d show the prediction of SOE using US06 drive cycle at four different temperature ranges, i.e. -10 °C, 0 °C, 10 °C, and 25 °C, respectively. Table 7 shows the values of RMSE and MAE for both the technique of machine learning, i.e. DNN and SVR.

#### 4.2 Case Study 2

In case 1 the performance of model was analysed on a new drive cycle which was not fed for the training. In this case, model was evaluated using the drive cycles at temperatures set other than which fed for training. In this case all the drive cycles were used for the training but only at three temperature values, i.e. -10, 0, and 25 °C. Drive cycles at 10 °C were kept for the testing of developed model accuracy. Three drive cycles at 10 °C, i.e. UDDS, LA92, and HWFET was used for the validation



**Fig. 5** Prediction of SOE using DNN and SVR technique on US06 drive cycle at following temperature range **a** at -10 °C temperature **b** at 0 °C temperature **c** at 10 °C temperature **d** at 25 °C temperature

Table 7         Evaluation results on the individual drive cycles	
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	DNN		SVR	
Testing Drive Cycle	RMSE	MAE	RMSE	MAE
US06 at -10°C	0.0148	0.0125	0.0602	0.0442
US06 at 0°C	0.0149	0.0118	0.0898	0.0672
US06 at 10°C	0.0178	0.0128	0.0287	0.0218
US06 at 25°C	0.012	0.010	0.0164	0.0122

Table 8 Evaluation results on the combination of drive cycles

,					
	DNN		SVR		
Training Drive Cycle	RMSE	MAE	RMSE	MAE	
UDDS at 10 °C	0.0154	0.0120	0.0207	0.0156	
LA92 at 10 °C	0.0189	0.0162	0.0257	0.0177	
HWFET at 10 °C	0.0118	0.0093	0.0339	0.0235	



**Fig. 6** Prediction of SOE using DNN and SVR technique on following drive cycle at 10 °C temperature **a** UDDS drive cycle **b** LA92 drive cycle **c** HWFET drive cycle

purpose. This same procedure was followed in both ML techniques. Figure 6 shows the predicted SOE behaviour for the different combinations of drive cycles in graphical way. In this case also, it was seen that the error in the prediction of SOE was high for SVR method in comparison to DNN. The RMSE and MAE error values are shown in Table 8 for both methods and it can be easily seen that DNN performance is significantly better as compared to SVR.

### 5 Conclusion

This work offers a unique contribution to the accurate estimation of SOE. In this study modified algorithm for DNN and SVR with optimization techniques are implemented. The effectiveness of DNN and SVR methods were compared for the prediction of SOE. Instead of using any random hyperparameters, this study proposed the approach to optimize the hyperparameters and then used the optimized values of hyperparameters in the subsequent prediction. During the optimization, it was noticed that the error in the prediction reduces as the hidden layers and number of neurons increases in the case of DNN but higher values of them make the model computationally less efficient. Similarly in the case of SVR technique, if the value of regularization parameter (C) and gamma  $(\gamma)$  increases then the training time increases significantly, this makes the model less time efficient. This work shows the unique method for the optimization of hyperparameters of both techniques, i.e. DNN and SVR. The prediction result of this work shows that the DNN model is more efficient in the prediction of SOE as compared to SVR. The results obtained in the different case studies show that the SOE prediction by DNN model at a drive cycle different from the one used in training is quite good. In the same way, DNN model is able to predict the SOE at different thermal ranges other than those used in training. The results of the case studies evidently prove that the prediction from DNN is far better than SVR. This work also suggests that the conventional regression model used for the estimation of SOE can be upgraded using the DNN model in the future. This work highlights the prevailing challenges in the industry and proposes the potential recommendation for BMS development and SOE estimation in next-generation EV applications.

### Appendix A

# A.1 Modified DNN Algorithm With the Optimization Technique

- 1 Import dataset // Import the dataset required for the training
- 2 X = ['Current', 'Voltage', 'Average Current', 'Average Voltage'] // Input parameters
- 3 Y = ['SOE'] // Output parameters
- 4 Define range of values of hidden layers // range of values, example- 2, 3, 4.....
- 5 Define range of values of neurons // range of neurons values, example-1, 2, 4 .....
- 6 // Initializing optimization loop for hidden layers and neurons
- 7 For i in range of hidden layers values // loop for the range of hidden layers values
- 8 For j in range of neurons values // loop for the range of neurons values
- 9 find RMSE // rmse for the combination of hidden layer and neurons values
- 10 Hidden layers and neurons = (L, N) // Optimum values on the basis of minimum of rmse
- 11 Import model // import the model of DNN
- 12 From layers import Dense // import dense layer that will be add in model
- 13 model. add(Dense(units = N, input\_dim = no. of input parameters, activation = 'relu')) // add the dense layer, where N is optimum value of neuron, input\_dim is the no, of input parameters and activation function is ReLU. More layers can be add in the similar manner
- 14 model.add(Dense(units = 1, activation = 'linear')) // Output layer with Linear activation
- 15 model.compile(optimizer = Adam(learning\_rate = 0.01), loss = 'MSE', metrics = ['mae'])
   // Compile the model with Adam optimizer, learning rate fixed to be 0.01, loss function be
   MSE and metrics is MAE
- 16 model.fit(X, Y, epochs = 10000) // fitting of model where epochs is the no. of iterations
- 17 Import testing dataset // import the dataset for the testing of performance of developed model
- 18 Define X\_test and Y\_test // define the X\_test and Y\_test in the similar manner as done for training X and Y
- 19 Y\_predict = model.predict(X\_test) // Calculate the predicted output
- 20 RMSE(Y\_predict, Y\_test) // Finally calculate the rmse for predicted and actual output

# A.2 Modified $\upsilon$ —SVR Algorithm With the Optimization Technique

- 1 Import dataset // Import the dataset required for the training
- 2 X = ['Current', 'Voltage', 'Average Current', 'Average Voltage'] // Input parameters
- 3 Y = ['SOE'] // Output parameters
- 4 Define range of values of C // range of values selected were 0.01, 0.1, 1, 10, 100
- 5 Define range of values of  $\gamma$  // range of  $\gamma$  values selected 0.1, 1, 10
- 6 // Initializing optimization loop for hidden layers and neurons
- 7 For i in range of  $\gamma$  values // loop for the range of  $\gamma$  values
- 8 For j in range of C values // loop for the range of C values
- 9 find RMSE // rmse for the combination of C and  $\gamma$  values
- 10 C and  $\gamma = (c, gp) // Optimum values on the basis of minimum of rmse$
- 11 Import model // import the v-SVR model of SVM
- 12 model = NuSVR(kernel = 'rbf', C = c, gamma = gp, nu = 0.5) // fitting of model
- 13 Import testing dataset // import the dataset for the testing of performance of developed model
- 14 Define X\_test and Y\_test // define the X\_test and Y\_test in the similar manner as done for training X and Y
- 15 Y\_predict = model.predict(X\_test) // Calculate the predicted output
- 16 RMSE(Y\_predict, Y\_test) // Finally calculate the rmse for predicted and actual output

# **Appendix B**



# **B.1 US06 Drive Cycle Velocity Profile**

**B.2 UDDS Drive Cycle Velocity Profile** 



## **B.3 HWFET Drive Cycle Velocity Profile**



**B.4 LA92 Drive Cycle Velocity Profile** 



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