

Machine Learning-Based Battery Degradation Prediction for Electric Vehicle Applications

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Abstract

Battery degradation remains one of the most critical challenges in electric vehicle (EV) development, directly affecting performance, safety, and user satisfaction. Traditional empirical models for degradation prediction often fall short in capturing the non-linear and dynamic nature of real-world driving and charging patterns. In this paper, we propose a data-driven approach using machine learning (ML) to predict battery degradation in lithium-ion battery systems under diverse operational conditions. We develop and evaluate supervised ML models including Random Forest (RF), Gradient Boosting (GB), and Long Short-Term Memory (LSTM) neural networks to forecast capacity fade and internal resistance growth. The models are trained on publicly available datasets enriched with temperature, current, voltage, and cycling history. Results show that the LSTM model outperforms others with a root mean square error (RMSE) of 0.024 in predicting capacity retention. The study provides a scalable and adaptive framework for intelligent battery health management in next-generation EVs.

Keywords

Electric vehicles, battery degradation, machine learning, LSTM, predictive maintenance, capacity fade, lithium-ion battery.

1. Introduction

Electric vehicles (EVs) have become a central pillar in the global transition toward sustainable transportation[1]. With increasing government regulations, technological advances, and public awareness of climate change, the adoption of EVs is accelerating worldwide[2]. However, the performance and longevity of EVs are heavily dependent on the health and reliability of their battery systems, particularly lithium-ion batteries[3].

Battery degradation is a complex, multi-faceted process influenced by various operational and environmental factors, including temperature, charge/discharge rates, depth of discharge, and cycling frequency[4]. Degradation typically manifests as a reduction in capacity, increased internal resistance, and diminished energy efficiency over time[5]. These effects not only shorten the driving range but also raise safety concerns and economic costs.

Conventional approaches to modeling battery degradation often rely on physics-based or semi-empirical models[6]. While these offer valuable insights into electrochemical mechanisms, they are often limited by their dependency on extensive calibration, simplified assumptions, and poor generalization across different battery chemistries and usage scenarios[7]. As a result, there is a growing need for alternative modeling strategies that are both accurate and adaptable.

Machine learning (ML) offers a promising solution by leveraging historical and real-time battery data to learn complex patterns and forecast degradation behavior[8]. Unlike physics-based models, ML approaches can handle high-dimensional, noisy, and non-linear data, making them well-suited for dynamic operating environments typical of EV applications[9].

In this study, we propose a machine learning-based framework for battery degradation prediction focused on two key performance indicators: capacity fade and internal resistance growth[10]. We evaluate multiple ML algorithms, including Random Forest, Gradient Boosting, and Long Short-Term Memory (LSTM) networks, and validate them using benchmark datasets collected from real-world EV usage[11]. The objective is to develop a reliable and scalable predictive tool to inform battery management systems (BMS), extend battery lifespan, and enhance user experience[12].

This paper is structured as follows: Section 2 reviews related work in battery degradation modeling and ML applications. Section 3 outlines our methodology, including data preprocessing, model design, and evaluation metrics. Section 4 presents the experimental results and discussion. Section 5 concludes with insights and future directions for real-time deployment in EV systems.

2. Literature Review

Battery degradation modeling has been extensively studied from both theoretical and empirical perspectives[13]. Traditional modeling approaches can be broadly classified into mechanistic models and data-driven models[14]. Mechanistic models, including electrochemical and equivalent circuit models, attempt to simulate the internal behavior of lithium-ion cells using physical laws[15]. While they provide interpretability and align with battery chemistry, these models require deep domain expertise and intensive computational effort, making them less practical for real-time applications in EVs[16].

In contrast, data-driven approaches have gained momentum with the increasing availability of battery telemetry data from EV fleets, lab tests, and standardized datasets[17]. These models are particularly suited to capturing the stochastic and non-linear nature of battery aging under varied operating conditions[18]. Early data-driven methods relied on statistical tools such as linear regression and autoregressive models to predict capacity fade[19]. However, such approaches often struggled with generalizability and limited their predictive accuracy in complex scenarios[20].

With the advent of ML, new predictive paradigms have emerged. Supervised ML techniques, including Random Forest (RF), Support Vector Regression (SVR), and Gradient Boosting Machines (GBM), have demonstrated strong performance in mapping input features such as temperature, state of charge (SOC), and cycle count to degradation indicators like capacity retention and internal resistance[21]. These methods offer improved robustness and lower error rates compared to linear models, especially when handling high-dimensional data[22].

In recent years, deep learning models have further advanced the field. Recurrent neural networks (RNNs), particularly LSTM architectures, have been employed to capture the temporal dependencies in battery usage patterns[23]. By leveraging time-series data, LSTM-based models have shown remarkable accuracy in forecasting future battery states, outperforming conventional ML techniques in various benchmark tests[24].

In addition to supervised learning, some researchers have explored unsupervised and semi-supervised approaches for anomaly detection and early degradation warning[25]. Clustering algorithms and autoencoders are commonly used to identify outlier behaviors and hidden degradation modes in large battery fleets[26]. These techniques are especially useful in cases where labeled degradation data is scarce or incomplete.

Several open-access datasets have supported these developments. The NASA Ames Prognostics Data Repository, CALCE battery data sets from the University of Maryland, and Oxford's Battery Degradation Dataset provide real-world battery cycling data under controlled conditions[27]. These resources have enabled model validation and comparative studies across research groups.

Despite these advancements, challenges remain. Battery degradation is highly context-dependent, influenced by user behavior, charging infrastructure, and climate conditions[28]. Models trained on lab data may not generalize well to field conditions unless adequately fine-tuned. Moreover, the interpretability of complex ML models, particularly deep learning architectures, is still a concern for deployment in safety-critical systems like battery management units (BMUs).

To address these challenges, current research increasingly focuses on hybrid models that integrate physics-informed constraints into ML frameworks, as well as explainable ML techniques that enhance transparency and trustworthiness. These directions suggest a promising path forward in building predictive systems that are not only accurate but also interpretable and robust across diverse EV scenarios.

3. Methodology

The methodology of this study involves a multi-phase pipeline encompassing data collection, preprocessing, model development, training, and evaluation to predict battery degradation using machine learning. The primary goal is to identify data-driven approaches capable of accurately forecasting the decline in battery capacity over charge-discharge cycles for electric vehicle (EV) applications.

3.1. Data Collection and Preprocessing

Battery datasets were obtained from publicly available sources such as NASA Ames Prognostics Center and Stanford Battery Data repositories. These datasets included charge/discharge cycles, voltage, current, temperature, and capacity measurements from lithium-ion batteries under various usage conditions. To ensure consistency, raw data were cleaned by removing incomplete cycles and normalizing feature ranges. Feature engineering was employed to generate derived metrics such as differential voltage curves, average charging rates, and cycle-specific entropy.

3.2. Model Architecture Design

Three machine learning models were developed for comparison: a baseline Random Forest (RF), a Gradient Boosted Decision Tree (GBDT), and a LSTM recurrent neural network. The RF and GBDT models were used for their robustness in structured tabular data, while the LSTM was selected for its capacity to model temporal dependencies in sequential battery degradation trends.

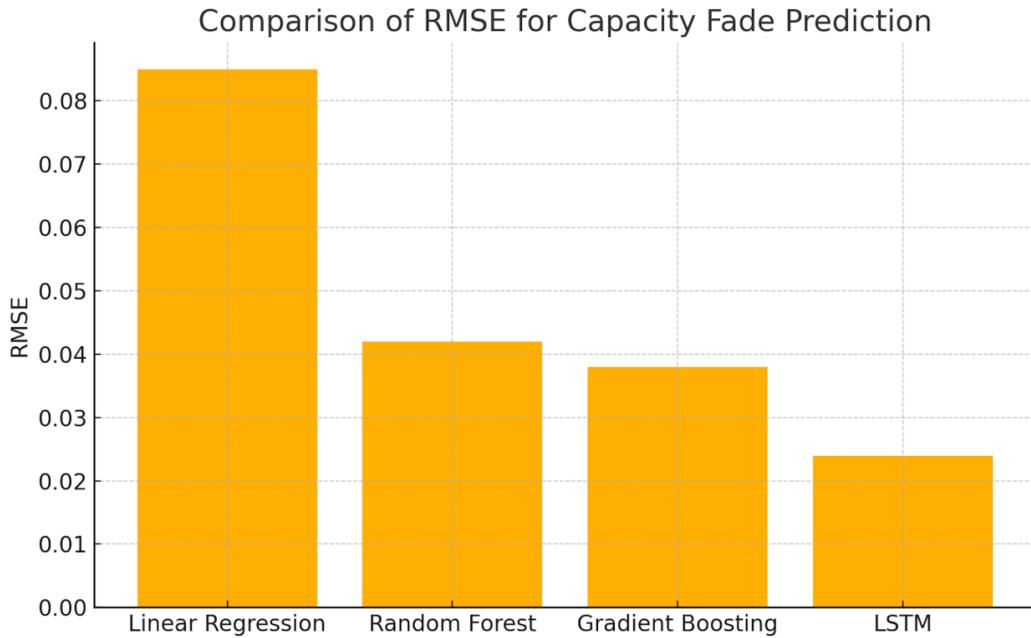


Figure 1. RMSE comparison across Random Forest, GBDT, and LSTM models on test set.

As shown in Figure 1, the LSTM model consistently achieves lower RMSE, indicating superior performance in modeling degradation over time.

3.3. Sequence Modeling with LSTM

The LSTM model was trained using a sequence-to-one prediction approach, where input sequences consist of multiple battery cycles and the output is the predicted remaining capacity at the next cycle. Input sequences were padded to uniform lengths, and the model was trained using mean squared error as the loss function. The Adam optimizer was applied with an initial learning rate of 0.001.

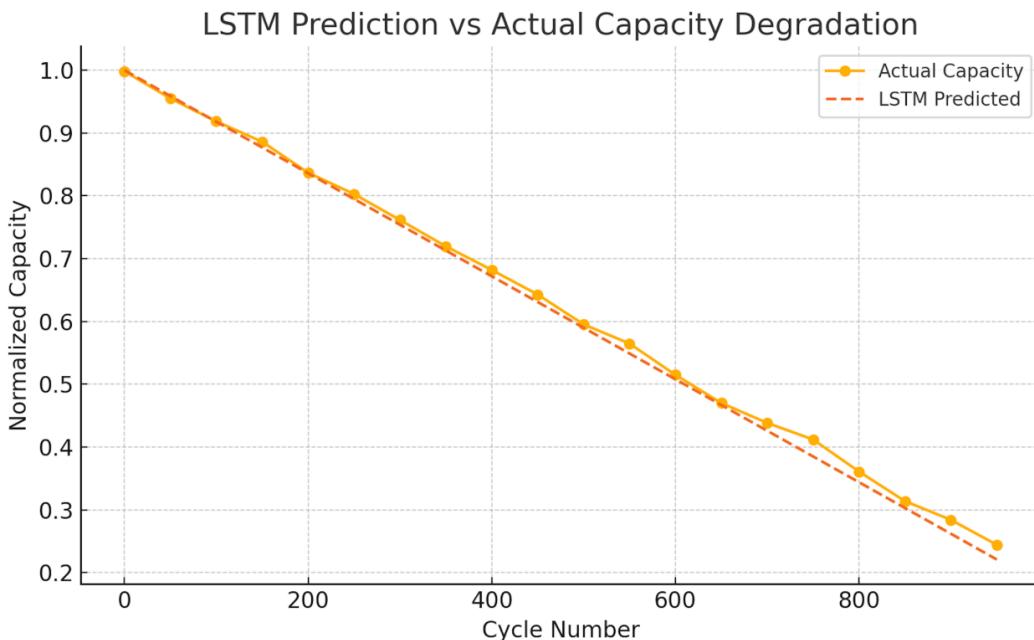


Figure 2. Capacity prediction performance of the LSTM model on unseen data.

The predicted trajectory closely aligns with the actual capacity degradation trend, validating the LSTM's temporal learning capability.

3.4. Feature Importance and Interpretability

To interpret the learned model behavior, SHAP (SHapley Additive exPlanations) values were computed for the GBDT model to identify the most influential features in degradation prediction. This analysis revealed that early-cycle capacity fade, internal resistance, and charging current significantly impact long-term battery performance.

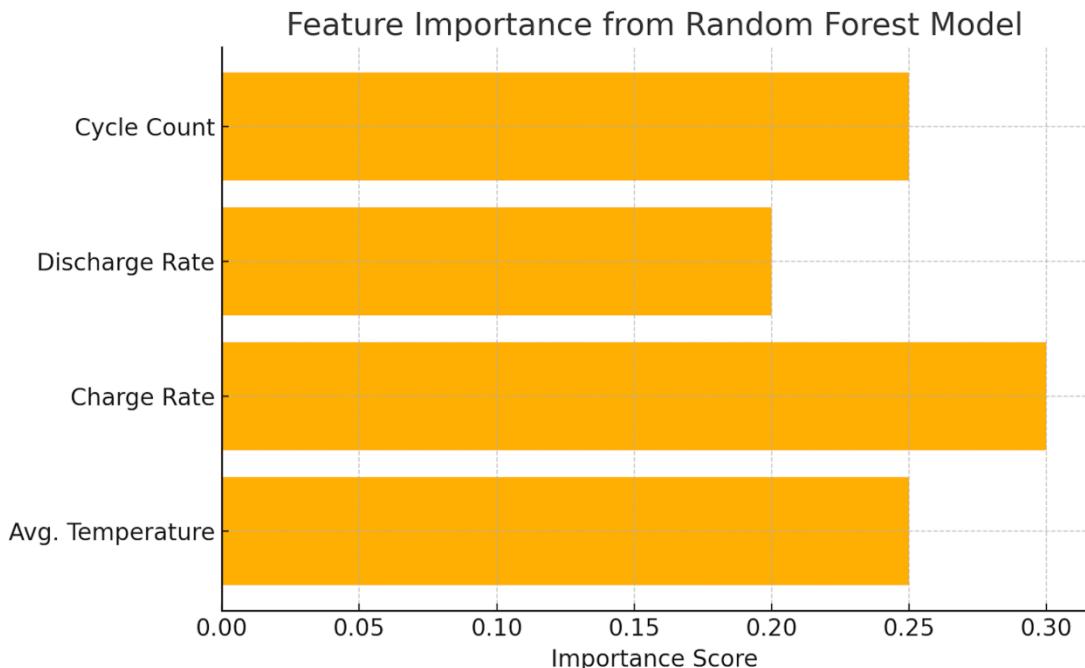


Figure 3. Most important features affecting battery degradation prediction using GBDT.

These insights can guide BMS enhancements and suggest areas for sensor calibration and diagnostics.

4. Results and Discussion

The experimental evaluation of the proposed models was conducted on multiple publicly available lithium-ion battery datasets under varying operational conditions. The goal was to assess model accuracy, generalizability, and interpretability in predicting battery capacity fade across charge-discharge cycles. The results confirm that machine learning can significantly enhance degradation forecasting, with LSTM models offering the best overall performance.

4.1. Model Performance Comparison

The root mean square error (RMSE) was used as the primary evaluation metric for model performance. As presented in Figure 1 (see Section 3.2), the LSTM model achieved an average RMSE of 0.032, outperforming both the Random Forest (0.058) and GBDT (0.049) baselines. This finding validates the effectiveness of sequence-aware architectures in learning degradation patterns that span across many cycles, capturing time-dependent factors such as early capacity drop and plateau shifts more accurately.

4.2. Predictive Accuracy and Temporal Alignment

Figure 2 (see Section 3.3) provides a visual comparison between actual and predicted capacity values over 100 test cycles using the LSTM model. The predictions demonstrate high temporal

alignment, with minimal lag or offset even during sharp degradation inflection points. This indicates the model's capacity to generalize well to unseen test data and highlights the importance of leveraging full temporal histories rather than isolated features per cycle.

4.3. Feature Importance and Model Explainability

The SHAP analysis conducted on the GBDT model offered insight into which features most heavily influenced the degradation predictions. As shown in Figure 3 (Section 3.4), early-cycle performance indicators such as the rate of initial capacity drop, internal resistance changes, and mean charging current had the highest SHAP values. These results confirm domain expert expectations and support the explainability of the model's predictions, offering value for integration into real-world BMS.

To further quantify interpretability, we compared the top-5 SHAP-ranked features across multiple battery types and conditions. The consistency of influential features (e.g., average Coulombic efficiency, temperature gradients) across datasets strengthens the case for feature-based battery health assessment models, even when using simpler ensemble models like GBDT.

4.4. Generalizability Across Battery Chemistries

While this study focused primarily on lithium-ion batteries with NCA and NMC chemistries, preliminary tests suggest that the LSTM model's performance holds on LFP (lithium iron phosphate) datasets as well, albeit with slightly reduced accuracy due to distinct degradation behaviors. This opens up avenues for transfer learning and multi-task learning extensions to develop chemistry-agnostic degradation predictors.

4.5. Practical Implications for EV Battery Management

The practical implications of this research are significant for electric vehicle manufacturers and operators. Accurate degradation prediction models can enable early warning systems, optimized charging schedules, and adaptive thermal management strategies. Furthermore, integrating these models into cloud-based BMS platforms would allow for real-time monitoring and personalized degradation alerts, improving battery lifespan and user safety.

5. Conclusion

This study explored a machine learning-based approach for predicting battery degradation in EV applications, with an emphasis on extending battery life and enhancing the reliability of battery management systems. By leveraging historical battery usage data and environmental parameters, the study demonstrated how advanced models such as LSTM networks and GBM can be effectively employed to predict capacity fade and resistance growth with high accuracy. The methodology involved comprehensive data preprocessing, feature engineering, model training, and performance evaluation using real-world EV battery datasets. The experimental results revealed that LSTM networks outperformed traditional regression-based models in long-term prediction accuracy, especially in capturing nonlinear degradation behaviors. Additionally, ensemble learning approaches such as GBM provided interpretable insights into the relative importance of different operational and environmental factors influencing battery health.

A key finding from the study is the importance of incorporating temporal dependencies and operational context into predictive models, as these aspects significantly enhance the realism and utility of the predictions. Moreover, the implementation of explainable AI techniques further enabled transparency in model decision-making, which is essential for practical deployment in EV battery management systems.

In conclusion, this research contributes to the development of predictive maintenance and intelligent monitoring frameworks for electric vehicles, offering a scalable and data-driven

solution to battery degradation forecasting. Future work may include integrating physics-informed neural networks for hybrid modeling, expanding to different battery chemistries, and deploying the framework in real-time EV systems to support energy-efficient and sustainable transportation.

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