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Hybrid ML-Based Cutting Temperature Prediction in Hard Milling Under Sustainable Lubrication

Balasuadhakar Arumugam [1], Thirumalai Kumaran Sundaresan [2,*] and Saood Ali [3,*]

[1] Department of Mechanical Engineering, Sona College of Technology, Salem 636005, Tamil Nadu, India
[2] Department of Mechanical Engineering, PSG Institute of Technology and Applied Research, Coimbatore 641062, Tamil Nadu, India
[3] School of Mechanical Engineering, Yeungnam University, 280 Daehak-Ro, Gyeongsan 38541, Gyeongbuk, Republic of Korea
* Correspondence: thirumalaikumaran@yahoo.com (T.K.S.); saoodali@ynu.ac.kr (S.A.)

Abstract

The field of hard milling has recently witnessed growing interest in environmentally sustainable machining practices. Among these, Minimum Quantity Lubrication (MQL) has emerged as an effective strategy, offering not only reduced environmental impact but also economic benefits and enhanced cooling performance compared to conventional flood cooling methods. In hard milling operations, cutting temperature is a critical factor that significantly influences the quality of the finished component. Proper control of this parameter is essential for producing high-precision workpieces, yet measuring cutting temperature is often complex, time-consuming, and costly. These challenges can be effectively addressed by predicting cutting temperature using advanced Machine Learning (ML) models, which offer a faster and more efficient alternative to direct measurement. In this context, the present study investigates and compares the performance of Conventional Minimum Quantity Lubrication (CMQL) and Graphene-Enhanced MQL (GEMQL), with sesame oil serving as the base fluid, in terms of their effect on cutting temperature. The experiments are structured using a Taguchi L36 orthogonal array, with key variables including cutting speed, feed rate, MQL jet pressure, and the type of cooling applied. Additionally, the study explores the predictive capabilities of various advanced ML models, including Decision Tree, XGBoost Regressor, K-Nearest Neighbor, Random Forest Regressor, and CatBoost Regressor, along with a Hybrid Stacking Machine Learning Model (HSMLM) for estimating cutting temperature. The results demonstrate that the GEMQL setup reduced cutting temperature by 36.8% compared to the CMQL environment. Among all the ML models tested, HSMLM exhibited superior predictive performance, achieving the best evaluation metrics with a mean absolute error of 3.15, root mean squared error (RMSE) of 5.3, mean absolute percentage error of 3.9, coefficient of determination (R^2) of 0.91, and an overall accuracy of 96%.

Keywords: minimum quantity lubrication; machining; cutting temperature; machine learning

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1. Introduction

AISI H11 hot work tool steel is widely used for machining tools, dies, hot shear blades, and aerospace components because of its high hardness, toughness, wear resistance, and reliable performance under high stress and elevated temperatures [1]. Milling is a versatile machining process that produces flat, helical, and curved surfaces. The main challenges in hard milling are significantly reduced tool life and compromised surface quality [2]. In

hard milling, lubrication is vital for limiting friction and dissipating heat at the contact zone between the tool and the workpiece, which helps maintain tool performance and improves the machined surface [3]. Minimum Quantity Lubrication (MQL) technique enhances sustainability in the milling process by using minimal lubricant, thereby cutting down fluid waste and disposal issues. It further contributes to eco-friendly machining through reduced resource usage and lower environmental footprint [4]. Due to their renewable origin, biodegradability, and eco-friendly characteristics, vegetable oils serve as effective substitutes for mineral cutting fluids. Their application through MQL further promotes green and sustainable machining practices [5].

Studies reveal that nanoparticles enhance lubricant performance by reducing friction and wear. This is achieved through mechanisms like rolling, self-repair, polishing, and tribo-film formation [6]. Haq et al. studied copper-based nanofluid MQL (NMQL) in milling Inconel 718 and found notable improvements, with a 14.7% drop in cutting temperature (T_c), 20.1% reduction in surface roughness (R_a), and 13.3% decrease in power consumption [7]. Pal et al. studied the impact of sunflower oil-based MQL with Al_2O_3 nanoparticles on the machinability of AISI 321 steel. Their findings indicated substantial reductions in key parameters compared to the conventional flood cooling method, including a 44% decrease in cutting force (F_c), a 67% reduction in torque (T), a 56% reduction in R_a, and a 26% decrease in T_c [8]. Dong et al. investigated the application of MoS_2 particles in an emulsion-based MQL system during the machining of SKD11 tool steel and reported a 55% reduction in surface Ra compared to dry machining [9]. Hamdi et al. studied the influence of hexagonal boron nitride (hBN) and multi-walled carbon nanotube (MWCNT) nanofluids in an MQL system with vegetable base oil for machining AISI H11. Compared to dry and pure MQL, MWCNTs lowered T_c and R_a, while hBN reduced energy consumption (U_c) [10].

Sesame oil, derived from sesame seeds, is rich in antioxidants and fatty acids, offering good lubrication and eco-friendly properties. Its biodegradability, availability, and low cost make it a sustainable option for multiple uses [11]. Padmini et al. studied MQL machining of AISI 1040 using sesame and coconut oils with MoS_2 additives, finding that sesame oil with nano MoS_2 gave the best results, with major reductions in tool wear (VB), R_a and F_c compared to dry cutting [12]. Furthermore, Ni et al. investigated the use of Fe_3O_4, Al_2O_3, and carbon nanoparticles dispersed in sesame oil for broaching AISI 1045 steel. Their results indicated that the carbon-based nanofluid exhibited higher viscosity, minimized friction, reduced F_c, lowered vibration, and improved surface finish compared to the others [13]. Graphene nanoplatelets (Gnps), with their two-dimensional layered structure, offer excellent wear resistance and friction-reducing behavior, making them effective micro- and nano-scale lubricants as well as additives in liquid lubricants [14]. Their unique atomic arrangement of carbon atoms in a honeycomb lattice also provides outstanding thermal conductivity. The strong carbon-carbon bonds allow heat to transfer rapidly through lattice vibrations (phonons) with minimal resistance, enhancing lubrication performance in both thermal and mechanical applications [15].

Touggui et al. explored the effectiveness of six distinct cooling conditions, including Gnps-enriched MQL, MoS_2-enriched MQL, MWCNT-enriched MQL, Gnps/MoS_2 hybrid MQL, MWCNT/MoS_2 hybrid MQL, along with dry cutting in the turning of AISI 304 stainless steel. The MQL system employed vegetable oil as its base fluid. The study revealed that Gnps-enriched MQL effectively reduced R_a, F_c, T_c, and VB, achieving greater improvements than other cooling environments, including hybrid nanofluids [16]. Paiva et al. studied the machining performance of SAE 52,100 steel using Gnps-enhanced synthetic and semi-synthetic fluids, in comparison with MQL and flood cooling. Their results revealed that the addition of Gnps led to a 29% reduction in R_a with synthetic oil and a 14% reduction in cutting energy [17]. Similarly, Pal examined the MQL drilling of AISI 321

stainless steel with vegetable-oil-based nanofluid containing Gnps and reported reductions in F_c (27.4%), T (64.9%), R_a (33.8%), and coefficient of friction (51.7%) compared to pure MQL, along with improved tool life [18]. Jiang et al. investigated the micro-bearing effect of Gnps when mixed with SiC in a hybrid nanofluid. Their findings demonstrated that this synergistic combination significantly enhanced the lubrication performance of the coolant [19]. Extending these findings, Li et al. investigated the milling of TC4 alloy using Gnps MQL in combination with a bionic micro-texture tool, which reduced F_c (40%), T_c (43.8%), and R_a (56.9%), while extending tool life by 112% compared to dry cutting [20].

Sustainable manufacturing is rapidly advancing through the adoption of machine learning (ML) and artificial intelligence, where these technologies are being applied to optimize processes, boost productivity, and reduce environmental impact [21]. Mia et al. proposed a predictive ML model based on the Artificial Neural Network (ANN) algorithm for estimating T_c during the turning of AISI 1060 steel under both dry and high cooling conditions. In this model, cutting speed (v_c) and feed rate (f) were considered as the primary input variables, and through comparative analysis, the ANN demonstrated superior accuracy over the Response Surface Model, as evidenced by its lower Mean Absolute Error (MAE) values [22]. Building upon this line of research, Gupta et al. conducted a comparative study employing both Response Surface Model and the Adaptive Neuro-Fuzzy Inference System to predict machining responses, including F_c, T_c, and R_a, under NMQL conditions. Their findings indicated that the Adaptive Neuro-Fuzzy Inference System provided enhanced predictive performance compared to the Response Surface Model, thereby highlighting the effectiveness of hybrid soft computing approaches in machining optimization [23].

Furthermore, Zhang et al. advanced this field by developing a Gaussian Process Regression model aimed at forecasting F_c, R_a, and tool life during turning operations. The model utilized v_c, f, and depth of cut as its input parameters, and it exhibited remarkable stability and accuracy, establishing itself as a reliable tool for predictive modeling in machining processes [24]. Jurkovic conducted a comparative study of Support Vector Regression (SVR), polynomial regression, and ANN for predicting R_a, F_c, and tool life in high-speed turning, and their result revealed that polynomial regression provided better prediction of R_a and F_c, while ANN resulted in the best prediction of tool life [25]. Another comparative study was conducted by Mahfouz et al. to predict R_a in the end milling of aluminum. The study evaluated the performance of SVR, K-Nearest Neighbor (KNN), and Decision Tree (DT) models, and the results showed that SVR attained the highest accuracy of 81.3% among the models [26]. Additionally, Sizemore developed an ANN model to predict surface roughness during the turning of germanium and compared its performance with DT, RF, AdaBoost, and SVR, demonstrating the competitive capability of ANN-based approaches in machining applications [27].

A pure ML algorithm relies on a single method or technique to address a problem from beginning to end. In contrast, a hybrid algorithm combines two or more approaches, often from different domains, to tackle a task or enhance the efficiency of a single method. Such integration allows hybrid techniques to exploit the advantages of each individual algorithm while compensating for their weaknesses [28]. Akbari et al. developed a physics-informed substructure approach combined with a Bayesian learning algorithm to predict milling machine tool stability in production settings, demonstrating improved accuracy even with limited or suboptimal training data [29]. In a related effort, Rahman et al. devised a physics-guided ML method for predicting the corrosion resistance of alloys, highlighting the growing integration of domain knowledge with data-driven models to enhance prediction reliability [30]. Furthermore, Kim et al. applied a Bayesian learning

method integrated with a Deep Multiscale Convolutional Neural Network, achieving significant improvements in forecasting *VB* during end milling operations [31].

Huange al. introduced two hybrid machine learning models, Particle Swarm Optimization integrated with SVR and Particle Swarm Optimization integrated with Least Squares SVR, to predict *VB* in milling. These models were compared with conventional SVR. The analysis showed that the hybrid model achieved the best predictive performance, with a mean absolute error of 5.1, a mean absolute percentage error of 0.018, and a root mean square error of 0.018 [32]. Furthermore, Karmi et al. developed three hybrid models, namely optimizer-based deep neural networks, genetic algorithm-based deep neural networks, and deep neural network with extended Kalman filters to predict F_c, R_a, and power consumption during the machining of EN GJL 250 grey cast iron. These hybrid methods were compared with traditional ML models such as SVM, DT, and Levenberg–Marquardt, and the findings showed that the deep neural network with extended Kalman filters achieved the highest predictive performance [33].

A comprehensive review of the existing literature has revealed research gaps in the area of T_c analysis under MQL conditions. Only a limited number of studies have examined this aspect. Additionally, investigations on sesame oil as a cutting fluid are scarce, and work involving Gnps-enriched sesame oil in MQL environments is even more limited. Furthermore, the application of advanced ML models for predicting T_c remains largely unexplored, while the utilization of hybrid ML algorithms in this context is highly limited. Consequently, the present investigation focuses on assessing the performance of Graphene-Enhanced MQL (GEMQL) relative to Conventional Minimum Quantity Lubrication (CMQL), employing sesame oil as the base fluid during the hard milling process. The cutting temperature is selected as the principal response parameter, with the study addressing the influence of f, V_c, lubrication environment, and MQL jet pressure (P_j). In addition, this work seeks to develop a predictive framework for T_c using a Hybrid Stacking Machine Learning Model (HSMLM). The predictive capability of the proposed HSMLM will be benchmarked against advanced ML models.

2. Materials and Methods

2.1. Data Acquisition

In this investigation, AISI H11 hot die steel was selected as the workpiece material. Its chemical composition consisted of C (0.40%), Si (0.98%), Mn (0.25%), P (0.021%), S (0.01%), Cr (4.88%), Mo (1.21%), V (0.42%), with Fe as the remaining balance. The milling trials were performed on an LMW CNC vertical machining center (Model: LV 45), India, as illustrated in Figure 1. A TiAlN-coated carbide end mill cutter was utilized, having four flutes, a diameter of 6 mm, a helix angle of 30°, a flute length of 20 mm, and an overall length of 64 mm [34]. The cutting process was carried out under two lubrication conditions, namely CMQL and GEMQL. A minimum quantity lubrication setup with a reservoir capacity of 3 L supplied the cutting fluid at a constant flow rate of 60 mL/h. Sesame oil was used as the base fluid for the GEMQL condition.

The graphene nanoplatelets (Gnps), procured from a commercial supplier, were incorporated as the nano-additive. Their specifications included a thickness of 5–10 nm, length of 5–10 µm, density of 0.279 g/cm^3, surface area ranging from 200–240 m^2/g, 4–10 layers, and a thermal conductivity of about 2000 W/m-K. Based on findings from earlier work, the nanofluid was prepared at a concentration of 0.8 wt.% Gnps [35]. The preparation followed a two-step approach: initially, the mixture underwent ultrasonic agitation for 30 min using a Lark ultrasonic bath; subsequently, it was homogenized with an iStir HP 550 magnetic stirrer at 600 rpm for 45 min [36].

Figure 1. Experimental setup.

Experimental trials were designed using the L36 orthogonal array. The parameters studied included V_c of 40, 50, and 60 m/min, f of 0.01, 0.02, and 0.03 mm/rev, and P_j of 2, 4, and 6 bar, which were selected considering tool manufacturer recommendations and evidence from the literature [37,38]. T_c was recorded by means of a K-type thermocouple linked to an NI DAQ 9212 data acquisition unit, operating at a sampling frequency of 96 samples per second per channel. To improve accuracy, the thermocouple tip was positioned at a uniform distance of 0.5 mm beneath the machined surface, with thermal paste (conductivity >1.93 W/m-K) applied in the gap to ensure efficient heat transfer [39]. Additionally, T_c data were validated using an FLIR E60 infrared thermal imaging camera, India, providing cross-verification for the thermocouple measurements [40].

2.2. Data Preprocessing and ML Model Generation

The dataset used in this study consists of 36 rows and 5 columns, and it is well-structured. It includes one dependent variable, T_c, and four independent variables: cooling condition, V_c, f, and P_j. Data preprocessing and ML models development were performed in Python (Python V3.12.12) using Google Colab. The dataset, stored in CSV format, was imported into the notebook, and its structure was verified by checking its shape, feature details, data types, and null values. The numerical variables V_c, f, and P_j, and T_c were isolated for further analysis. Their distributions were examined using matplotlib distribution plots, while violin plots were employed to detect the presence of outliers.

A correlation heat map was generated to evaluate the association between the dependent variable T_c and the independent variables V_c, f, and P_j. Furthermore, to address the issue of overfitting caused by the limited dataset and to enhance the accuracy of the ML models, Gaussian data augmentation was applied. A noise factor equal to 0.01 was used, the augmentation method employed the fixed random seed of 42 to ensure reproducibility, and the augmentation process was repeated ten times [41]. Subsequently, the independent and dependent variables were separated as x and y. To ensure uniformity, the dataset underwent normalization through the StandardScaler method. Finally, the dataset was divided into training and testing sets, allocating 80% for model training and 20% for evaluation [42].

The DT is a widely used ML method for both regression and classification tasks. It operates as a predictive model by recursively splitting the dataset into subsets according to the most significant features, ultimately forming a tree-like structure. The criterion commonly used for splitting nodes in decision trees is the Gini impurity. It is calculated using the following formula, where kq represents the proportion of samples in the node belonging to class q [43].

$$I_G(k) = 1 - \sum_{q=1}^{Q} k_q^2 \tag{1}$$

XGBoost is a powerful ML algorithm that combines DT with gradient boosting, providing efficient and precise results for both regression and classification applications [44].

$$\hat{y}_i^{(t)} = \sum_{k=1}^{t} f_{k(x_i)} = \hat{y}_i^{(t-1)} + f_t(x_i) \tag{2}$$

where $\hat{y}_i^{(t)}$ represents the final constructed tree model; $\hat{y}_i^{(t-1)}$ is the tree model generated previously; $f_t(x_i)$ denotes the recently produced tree model, and t is the overall count of base tree models. In XGBoost, tree depth and the number of trees are key parameters, and optimization involves finding a classifier that minimizes the chosen loss function, as shown in Equation (3).

$$obj^t = \sum_{i=1}^{t} L\left(y_i, \hat{y}_i^{(t)}\right) + \sum_{i=1}^{t} \Omega(f_i) \tag{3}$$

where y_i designates the actual value, $\hat{y}_i^{(t)}$ denotes the predicted value, the loss function is represented by $L\left(y_i, \hat{y}_i^{(t)}\right)$, and the regularization term is denoted as $\Omega(f_i)$.

The KNN method operates on the principle that data points with comparable characteristics are typically located close together within the feature space. To make predictions, KNN examines the 'k' closest instances from the training set and assigns the outcome based on the most common class among those neighbors. Its performance heavily depends on the distance measure used and the selection of an appropriate k value. In addition, the method often demands proper preprocessing, such as feature scaling or normalization, to ensure reliable results [45].

$$\text{Predicted value } \hat{y} = \frac{1}{k}\sum_{i=1}^{k} y_i \tag{4}$$

The Random Forest Regressor (RFR) is an ensemble-based approach capable of handling both regression and classification tasks. It constructs multiple decision trees by repeatedly splitting the dataset into more uniform nodes according to predictor variables. The collective prediction is then derived by averaging the outputs of all trees, thereby enhancing model accuracy [46].

$$D = \{(x_i, y_i)\}_{i=1}^{N} \tag{5}$$

$$D_n = \{(x_i, y_i)\}_{i \in n} \tag{6}$$

$$split(x, \varnothing, \theta) = \arg\min_{\varnothing,\theta}\left(\frac{N_l}{N_n} Var(D_l) + \frac{N_r}{N_n} Var(D_r)\right) \tag{7}$$

$$Var(D) = \frac{1}{|D|}\sum_{(x_i,y_i)\in D}(y_i - \overline{y})^2 \tag{8}$$

$$\hat{y}_m(x) = \frac{1}{\left|D_{leaf}\right|}\sum_{(x_i,y_i)\in D_{leaf}} y_i \tag{9}$$

$$\hat{y}(x) = \frac{1}{m}\sum_{m=1}^{M} \hat{y}_m(x) \tag{10}$$

$$\varnothing_j = \frac{1}{M}\sum_{m=1}^{M} \sum_{node\ split\ on\ \varnothing_j} \left(Var(D_n) - \frac{N_l}{N_n} Var(D_l) - \frac{N_r}{N_n} Var(D_r)\right) \tag{11}$$

$$\hat{y}_{OOB,i} = \frac{1}{|D_{OOB,i}|}\sum_{T_m \ni i}(y_i - \hat{y}_{OOB,i})^2 \tag{12}$$

where D_n is the dataset contained in a node, N_n is the count of instances within the node, N_l denotes the count of instances in the left side child node, N_r is the count of instances in the right side child node, and D_l represents the dataset in the left child node, while D_r denotes the dataset in the right child node, D_{leaf} is the set of training instances in the leaf node, $\varnothing_j$ represents the feature importance, and OOB is out-of-bag data.

CatBoost is a gradient boosting algorithm that uses binary decision trees and stands out with ordered boosting, efficient modifications, and strong handling of target leakage, making it well-suited for small datasets. This technique, referred to as Greedy Target-based Statistics, operates by iteratively selecting the most optimal target-based statistics to enhance model accuracy [47]. CatBoost can be described as follows:

$$\frac{\sum_{j=1}^{p}\left[x_{j,k} = x_{i,k}\right]y_i}{\sum_{j=1}^{n}\left[x_{j,k} = x_{i,k}\right]} \tag{13}$$

Consider a dataset consisting of observations $D = \{X_i, Y_i\}\ i = 1, \ldots, n$, given a permutation $\sigma = (\sigma_1 \ldots, \sigma_n)$, $x_{\sigma_{p,k}}$ is substituted using the following relation, where p is a prior estimate and a denotes the weight assigned to this prior estimate.

$$\frac{\sum_{j=1}^{p-1}\left[x_{\sigma_{j,k}} = x_{\sigma_{p,k}}\right]y_{\sigma_j} + ap}{\sum_{j=1}^{p-1}\left[x_{\sigma_{j,k}} = x_{\sigma_{p,k}}\right] + a} \tag{14}$$

Hyperparameter tuning involves optimizing predefined parameters of a machine learning algorithm that are not learned from data, aiming to improve model performance and generalization. A common method is grid search with cross-validation, which systematically tests combinations of hyperparameters to identify the best-performing setup [48].

$$\Theta = \left\{(\theta_1, \theta_2, \ldots, \theta_p)\,\middle|\,\theta_i \epsilon \Theta_i\right\} \tag{15}$$

$$D = \bigcup_{K=1}^{K} D_k \tag{16}$$

$$D_{Train,k} = \frac{D}{D_k} \tag{17}$$

$$M_\theta^{(k)} = Train(M, D_{Train,k}, \theta) \tag{18}$$

$$P_\theta^{(k)} = \mathbb{E}_{(x,y)} \epsilon\, D_{val,k}\left[\mathcal{L}\left(M_\theta^{(k)}, (x,y)\right)\right] \tag{19}$$

$$\overline{P}_\theta = \frac{1}{k}\sum_{k=1}^{K} P_\theta^{(k)} \tag{20}$$

$$\sigma_\theta^2 = \frac{1}{k-1}\sum_{k=1}^{K}\left(P_\theta^{(k)} - \overline{P}_\theta\right)^2 \tag{21}$$

$$\theta^* = \underset{\theta \epsilon \Theta}{\operatorname{argmax}} \overline{P}_\theta \tag{22}$$

$$\theta^* = \underset{\theta \epsilon \Theta}{\operatorname{argmax}} \left(\overline{P}_\theta - \lambda \sigma_\theta\right) \tag{23}$$

where D represents the dataset and Θ denotes the hyperparameter space, and the dataset is divided into k folds for cross-validation, with D_k indicating the k^{th} fold. For each fold k, the training set is designated as $D_{Train,k}$ and the validation set as $D_{val,k}$. The model parameterized by hyperparameters θ for the k^{th} fold is expressed as $M_\theta^{(k)}$, and the performance is assessed using the loss function $\mathcal{L}$. The optimal hyperparameters, θ^*, are those that

maximize the average performance across all folds. Additionally, a trade-off parameter λ is used to balance mean performance and variance, ensuring robust model evaluation and selection.

Data augmentation (DA) is a strategy used to expand a limited dataset by artificially generating new samples. Among various DA approaches, Gaussian Data Augmentation (GDA) is considered one of the most efficient methods. It enhances the dataset by adding Gaussian noise to existing data points, thereby reducing data scarcity issues and improving the model's ability to generalize. Applying GDA reduces overfitting, enhances robustness, and boosts prediction accuracy. Gaussian noise is a random variation that adheres to a normal distribution, typically defined by a mean (α) of 0 and a standard deviation (β) of 1. Here, x represents a random variable drawn from this distribution [49]. The probability density function describing Gaussian noise can be expressed mathematically as follows:

$$f(x) = \frac{1}{\sqrt{2\pi\beta}} exp\left(-\frac{(x-\alpha)^2}{2\beta^2}\right) = \frac{1}{\sqrt{2\pi}} exp\left(-\frac{x^2}{2}\right) \quad (24)$$

The proposed Hybrid Stacking Model integrates the strengths of multiple machine learning algorithms to enhance predictive accuracy and robustness, as illustrated in Figure 2. In this framework, both categorical and continuous features are preprocessed using a column transformer, where polynomial interaction terms and standardization are applied to continuous inputs for improved representation. At the base level, diverse learners—including KNN, DT, RFR, XGBoost, and CatBoost—capture varying structural patterns and nonlinear relationships within the data. Their prediction outputs are subsequently combined through a meta-learner, which balances bias–variance trade-offs and mitigates overfitting effects through regularization mechanisms [50].

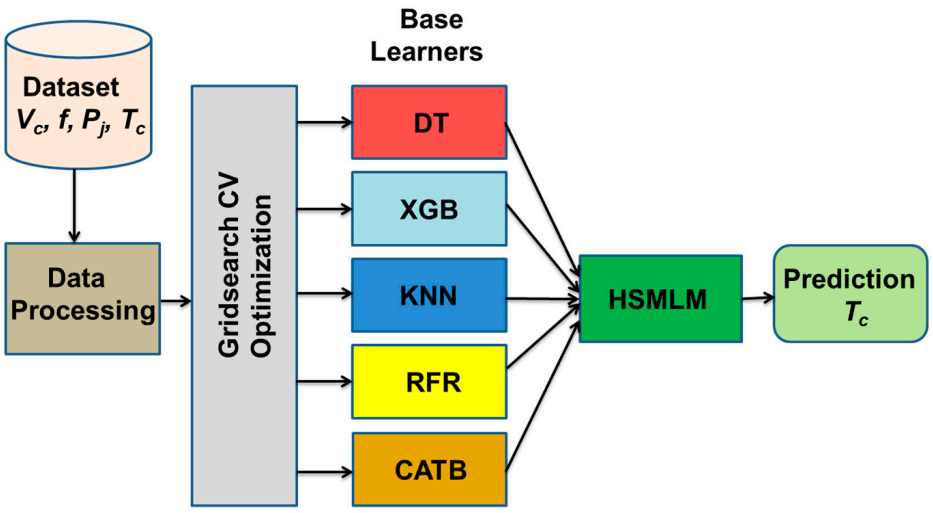

Figure 2. HSMLM Algorithm.

Additionally, a Transformed Target Regressor with Yeo–Johnson transformation is applied to stabilize variance and normalize the target distribution, ensuring smoother and more consistent model learning. Within this hybrid stacking architecture, layer-wise interaction occurs as the transformed feature vectors propagate through heterogeneous base learners, each capturing complementary predictive signals. The resulting prediction outputs are concatenated with the processed features to construct a richer meta-feature space for second-level learning, wherein the ElasticNet meta-learner assigns optimal weights by amplifying informative predictions while attenuating noisy contributions. During training, errors are iteratively minimized as the meta-learner refines residual patterns that are not captured

at the base layer. This hierarchical learning mechanism collectively improves robustness, enhances generalization, and delivers significantly improved predictive performance.

The continuous features are expanded with polynomial interaction terms and then standardized as expressed by:

$$\overline{x} = \varnothing(x) = \left[x^{(b)}, S\left(\varnothing_{poly}\left(x^{(c)}\right)\right)\right] \tag{25}$$

The meta-learner combines the base model predictions with the original preprocessed features using the ElasticNet regularization framework:

$$\min_{\beta,b} \frac{1}{2n}||z - (H\beta + b)||\binom{2}{2} + \alpha\left(\lambda||\beta||_1 + (1-\lambda)\frac{1}{2}||\beta||_2^2\right) \tag{26}$$

The final prediction is obtained by aggregating base learner outputs, applying the meta-learner, and then inverting the target transformation:

$$\hat{y}_* = T^{-1}(g([f_1(\phi(x_*)), \ldots, f_M(\phi(x_*)), \phi(x_*)])) \tag{27}$$

The binary features are denoted as $x^{(b)}$, while the continuous features are represented by $x^{(c)}$, including Speed, Feed, and MQL Pressure. The function $\varnothing_{poly}(.)$ refers to polynomial feature expansion, and $S(.)$ indicates standardization that ensures zero mean and unit variance. After these transformations, the final preprocessed feature vector is expressed as $\overline{x}$. The meta-feature matrix, which combines base learner predictions with preprocessed features, is represented as H. Within the meta-learner, β denotes the weight coefficients, b is the intercept term, α represents the overall regularization strength, and λ defines the penalties. For a new input sample, the notation x_* is used, and its preprocessed version is given $\phi(x_*)$.

Evaluation metrics are essential for assessing the effectiveness of ML models. This study uses five performance metrics: Coefficient of Determination (R^2), MAE, Mean Absolute Percentage Error (MAPE), Root Mean Square Error (RMSE), and Accuracy to identify the optimal model. In the equations, o_i represents the observed experimental values, while p_i represents the values predicted by the proposed ML models [51].

$$\text{MAE} = \frac{\sum_{i=1}^{n}|o_i - p_i|}{n} \tag{28}$$

$$\text{RMSE} = \sqrt{\frac{\sum_{i=1}^{n}(o_i - p_i)^2}{n}} \tag{29}$$

$$\text{R2} = 1 - \frac{\sum_{i=1}^{n}(o_i - p_i)^2}{\sum_{i=1}^{n}(o_i - \overline{o})^2} \tag{30}$$

$$\text{MAPE} = \frac{1}{n}\sum_{i=1}^{n}\frac{|o_i - p_i|}{o_i} \tag{31}$$

$$\text{Accuracy} = \left[1 - \left[\frac{1}{n}\sum_{i=1}^{n}\frac{|o_i - p_i|}{o_i}\right]\right] \times 100 \tag{32}$$

3. Results and Discussions

3.1. Influence of Lubrication Environment on Cutting Temperature

Examining T_c is essential, as it reveals how heat is generated and distributed in machining, influencing tool life and workpiece accuracy. This knowledge aids in selecting suitable cutting parameters, cooling methods, and tool materials to enhance efficiency and reliability [52]. Figure 3a depicts the effect of three different V_c at varying P_j on T_c.

It is observed that increasing the P_j generally reduces the T_c, while higher V_c leads to a substantial rise in T_c. This increase is attributed to the higher relative velocity between the tool and workpiece, which intensifies frictional heating at the cutting zone. Furthermore, elevated V_c intensifies plastic deformation in the workpiece material, thereby increasing thermal accumulation [53].

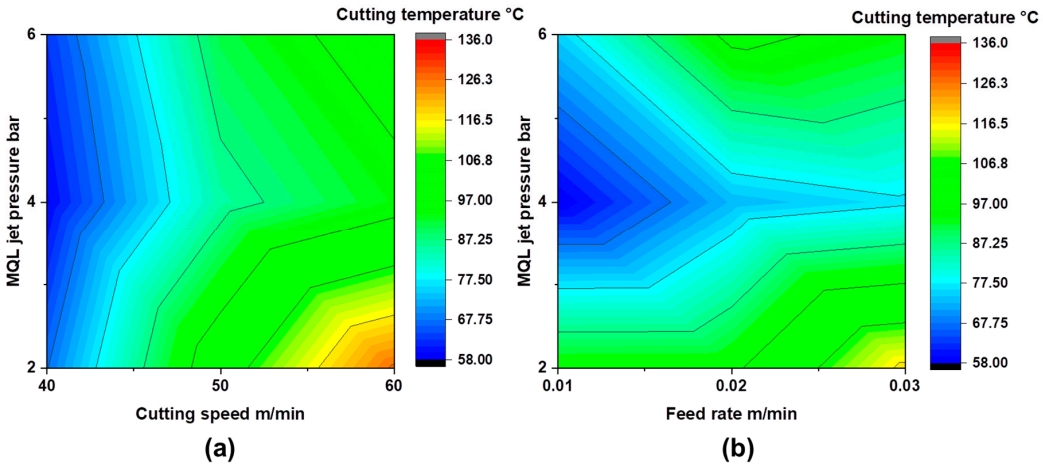

Figure 3. T_c at CMQL environment (**a**) V_c vs. P_j (**b**) f vs. P.

Higher P_j improves the cooling and lubrication effectiveness at the interface between the tool and workpiece. At low pressures, such as 2 bar, the mist does not penetrate efficiently into the cutting zone, allowing heat to accumulate and resulting in higher T_c [54]. Conversely, at excessively high pressures, such as 6 bar, the air pressure induces a spring-back effect, causing droplets to rebound from the cutting surface and reducing the effectiveness of lubrication. Thus, both very low and excessively high P_j can lead to elevated T_c. Optimal P_j provides sufficient penetration and heat dissipation, lowering the overall T_c. At low P_j, V_c strongly dominates the thermal response, causing a steep increase in T_c, whereas at optimal P_j, the rate of T_c rise is moderated due to enhanced cooling efficiency [55].

As shown in Figure 3b, T_c also increases with higher f; however, the underlying mechanism differs from that associated with variations in V_c, reflecting the distinct ways in which f influences the rise in T_c in the cutting zone. Higher f increases the uncut chip thickness, generating larger F_c and greater plastic deformation in the shear zone, which intensifies heat generation due to elevated shear strain and tool–chip friction [56]. P_j similarly regulates the thermal response under varying f. At low P_j, insufficient heat evacuation leads to thermal accumulation, whereas optimal P_j improves atomization and cooling, reducing the effective T_c. The interaction between f and P_j indicates that the adverse effect of higher f can be partially mitigated by increasing the P_j. Adequate P_j ensures that T_c remains within acceptable limits, even at higher f. This occurs because higher f generates more frictional heat, while an optimum P_j enhances coolant penetration, lubrication, and chip evacuation. Consequently, the additional heat is effectively dissipated, keeping T_c within controlled thermal boundaries even at elevated f levels [57]. From the contour plot analysis, the minimum T_c of 58 °C was achieved at a V_c of 40 m/min, f of 0.01 mm/rev, and P_j of 4 bar. Conversely, the maximum T_c of 136 °C occurred at 60 m/min V_c, 0.01 mm/rev f, and 2 bar P_j. Therefore, under the CMQL lubricating environment, the optimal cutting condition for minimizing T_c was identified as a V_c of 40 m/min, f of 0.01 mm/rev, and P_j of 4 bar.

The contour plot in Figure 4 demonstrates how T_c responds to variations in machining parameters under a GEMQL environment. Specifically, Figure 4a highlights the relationship

between V_c and P_j, showing their combined impact on the temperature generated during the cutting process. The GEMQL lubricating environment significantly reduces T_c across all machining conditions compared to the CMQL environment. This improvement is primarily attributed to the enhanced lubrication properties provided by Gnps, which improve the fluid's thermal stability, retainability in the cutting zone, and the formation of a protective tribo-film on the tool–workpiece interface [58]. At lower P_j, such as 2 bar in the CMQL setup, the cooling effect is insufficient; however, the presence of Gnps in GEMQL compensates for this limitation by promoting effective heat dissipation and reducing friction. Similarly, the spring-back effect observed at 6 bar MQL pressure under CMQL conditions is mitigated in GEMQL due to the rolling action of Gnps, which facilitates smoother tool–chip interaction and further decreases T_c in this high-pressure region [59].

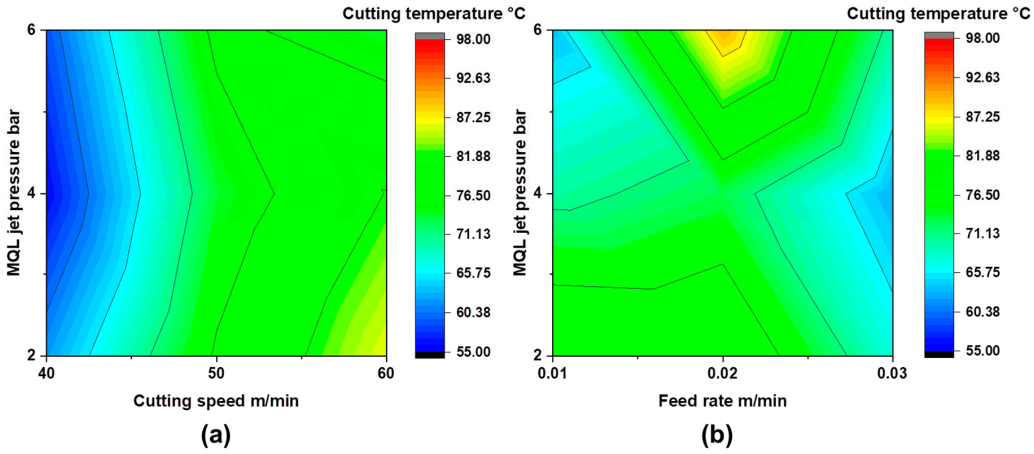

Figure 4. T_c at GEMQL environment (**a**) V_c vs. P_j (**b**) f vs. P.

As illustrated in Figure 4b, lower T_c values are achieved at both higher and lower feed rate ranges. The maximum T_c recorded was 98 °C at a V_c of 60 m/min, a f of 0.02 mm/rev, and 2 bar P_j. While this temperature is elevated under these conditions, it remains lower than the corresponding CMQL condition, which reached 106 °C. The maximum reduction in T_c of 36.8% was observed in the GEMQL environment compared to the CMQL environment at a V_c of 60 m/min, a f of 0.01 mm/rev, and a P_j of 2 bar. Similarly, Danush et al. reported a 31% reduction in T_c while milling Inconel 718 using a Gnps-based sunflower oil MQL system [60]. Analysis of the contour plots indicates that the optimal cutting conditions in the GEMQL environment are a V_c of 40 m/min, a f of 0.03 mm/rev, and a P_j of 4 bar.

3.2. Insight into Data Processing

Effective data processing plays a pivotal role in ML by converting raw data into informative representations that maximize a model's predictive capability. The distribution plot presented in Figure 5 illustrates the spread of the independent variables, namely V_c, f, and P_j. The V_c values considered in the dataset are 40 m/min, 50 m/min, and 60 m/min, while the f values are 0.01 mm/rev, 0.02 mm/rev, and 0.03 mm/rev. In a similar manner, the P_j values are 2 bar, 4 bar, and 6 bar. These independent variables are uniformly distributed, ensuring that the dataset provides an unbiased representation of the selected ranges.

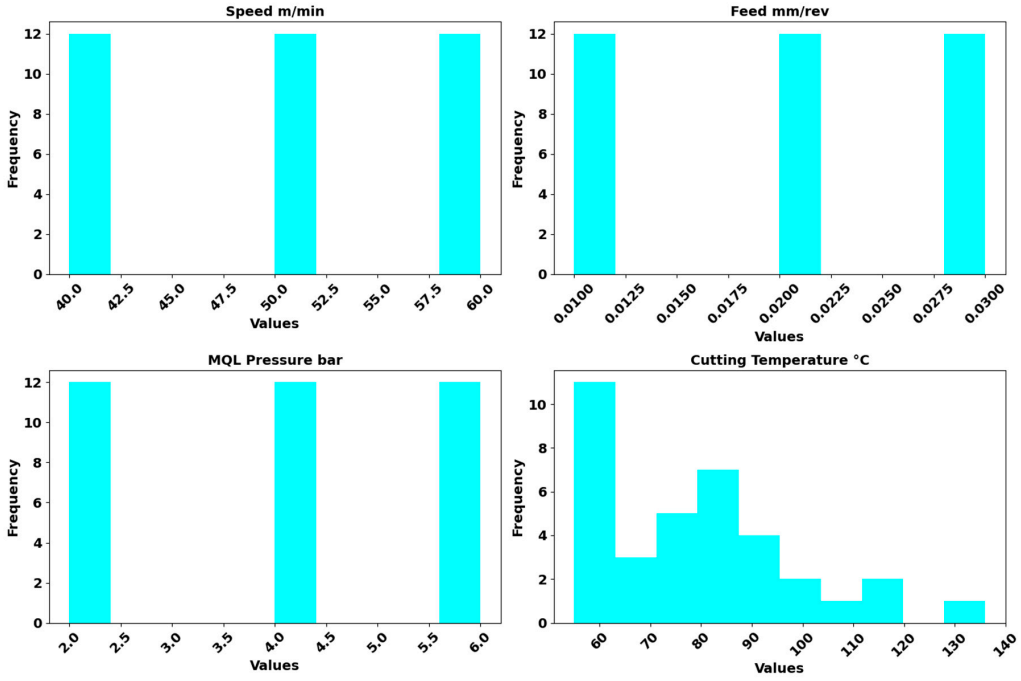

Figure 5. Distribution of independent and dependent variables.

In contrast, the dependent variable, T_c, exhibits noticeable skewness, indicating that its distribution is not symmetric. This skewness arises due to the presence of two distinct lubricating conditions, namely CMQL and GEMQL, which influence the temperature distribution differently under varying cutting parameters. To further assess the data quality, outlier analysis was performed using violin plots, as shown in Figure 6. The analysis revealed that no significant outliers were present in either the independent or dependent variables, thereby confirming the reliability of the experimental dataset. The absence of outliers is crucial, as extreme values can skew model training and lead to biased predictions, reducing overall accuracy.

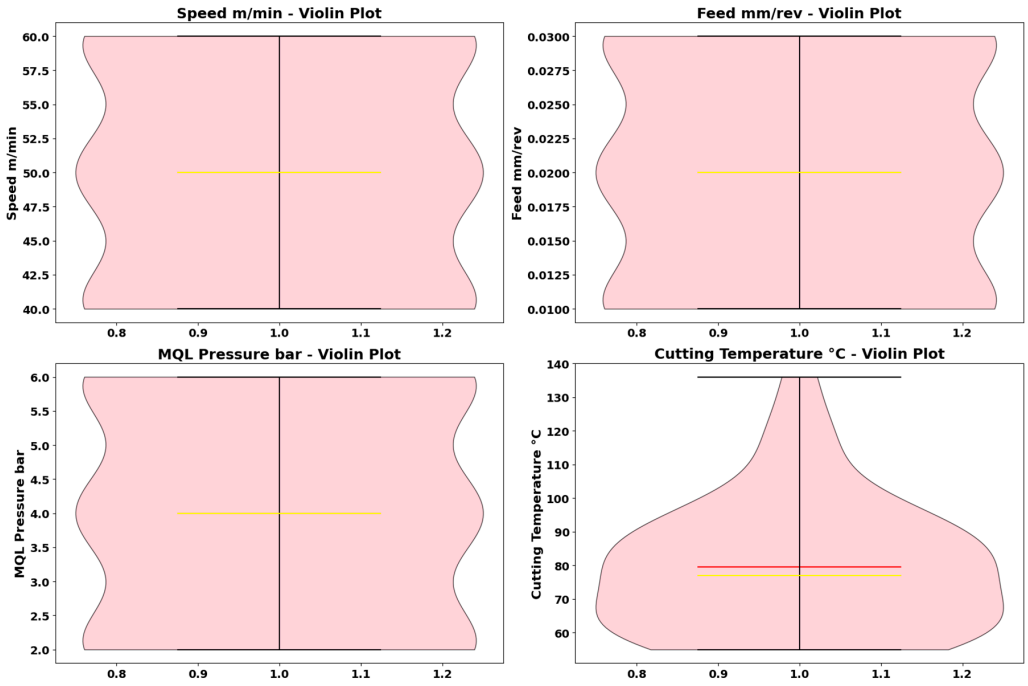

Figure 6. Violin plot for outlier detection.

The correlation analysis between the dependent variable (T_c) and the independent variables is illustrated in Figure 7, as examining these relationships helps identify influential features and improves model performance. The results demonstrate that T_c has the strongest positive correlation (0.7) with V_c, suggesting that an increase in V_c significantly elevates the T_c. A very weak positive correlation (0.033) was observed between T_c and f, implying that the effect of f on T_c is minimal. On the other hand, a moderate negative correlation (-0.17) was identified between T_c and P_j, indicating that higher P_j contributes to reducing the T_c to some extent.

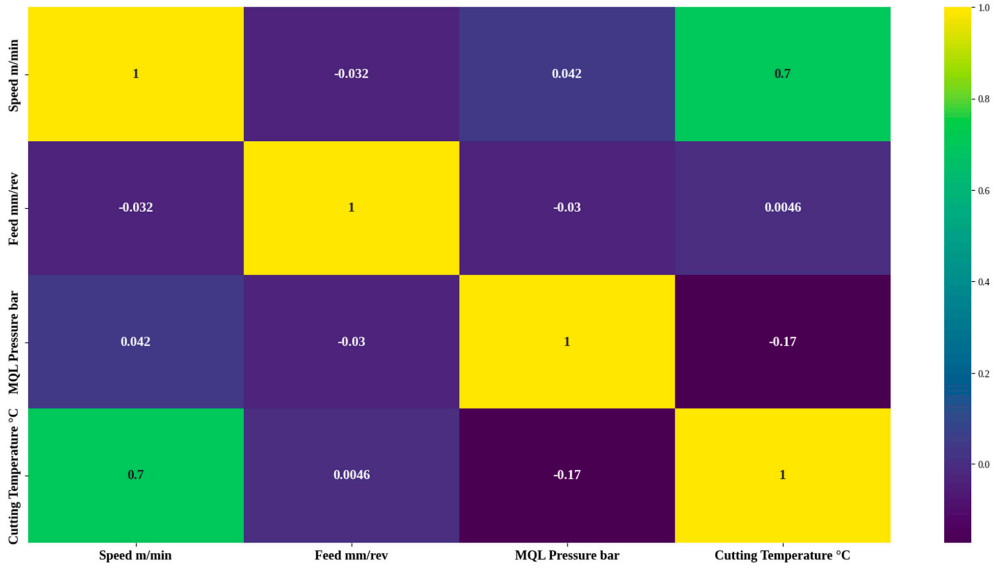

Figure 7. Correlation heat map.

3.3. Development and Comparison of ML Models

In this study, five advanced ML models, DT, XGBoost, KNN, RFR, and CatBoost, were initially developed as individual predictors and subsequently combined to create a hybrid stacked ML framework. In the DT model, the GridSearchCV tuning process considered three key parameters: maximum depth with values 16, 17, and 18; minimum samples per leaf with values 6, 7, and 8; and minimum samples per split with values 1, 2, and 4. The cross-validation was performed with three folds to ensure the reliability of the results. Maximum depth controls the overall complexity of the tree, while minimum samples per leaf regulates the minimum number of observations at the terminal nodes, preventing overly specific rules. Similarly, a minimum sample per split ensures that a node must have a sufficient number of samples before it can be divided further, thereby improving generalization. Based on this process, the optimal combination of hyperparameters was identified as max_depth = 17, min_samples_leaf = 8, and min_samples_split = 2. The performance of the tuned model was assessed using negative mean squared error (NMSE), which yielded a value of –91.48, demonstrating the effectiveness of parameter optimization in improving model accuracy.

The XGBoost model was tuned using five-fold cross-validation to identify the optimal hyperparameter. The values explored in the search included maximum depth with values [3, 5, 7, 10], number of estimators with values [50, 100, 200], learning rate with values [0.01, 0.05, 0.1, 0.2], subsample with values [0.7, 0.8, 1.0], and column sample by tree with values [0.7, 0.8, 1.0]. Based on this comprehensive search, the best parameters identified were a maximum depth of five, a learning rate of 0.1, a number of estimators was 200, a subsample value of 0.7, and a column sample by tree value of 1.0.

In the KNN model, the similarity between data points was measured using a distance metric, and the outcome was determined based on either equal weighting or distance-based weighting of the neighbors. The parameter "grid" explored included the number of neighbors with values [3, 5, 7, 9, 11], the weighting scheme with options "uniform" and "distance," and the distance metric with options "Euclidean" and "Manhattan." The grid search systematically evaluated all parameter combinations using two-fold validation, with performance measured through negative mean squared error. From this optimization process, the best parameters were identified as a distance metric of Manhattan, a number of neighbors equal to three, and a weighting scheme of distance. This configuration achieved an NMSE of -69.94.

The RFR model parameter "grid" was designed to explore different combinations of key hyperparameters. The parameter "number of estimators" was tested with values [50, 100, 200], which determines how many decision trees are included in the forest. The parameter "maximum depth" was set to [5, 10, 20] to control how deep each tree could grow, balancing model complexity with generalization. The parameter "minimum samples split with values" [2, 5, 10] specifies the least number of samples required to divide an internal node, while "minimum samples leaf with values" [1, 2, 4] determines the smallest number of samples allowed in a leaf node. These parameters play a key role in preventing overfitting and enhancing the model's stability. After the extensive search of GridSearchCV, the optimal configuration was determined to be a maximum depth of 20, 1 sample per leaf, 5 samples required to split a node, and 200 estimators. The optimized model achieved the best score of NMSE as -69.74.

The parameter "grid" for the CatBoost model was formulated to investigate various values of three key hyperparameters. The learning rate was tested with values [0.1, 0.2, 0.3], which controls the step size at each boosting iteration and balances the trade-off between speed of learning and overfitting. The depth was varied across [6, 8, 10], representing the maximum depth of the decision trees that form the base learners, with deeper trees allowing the model to capture more complexity. The iterations were set to [100, 200, 300], which determines the total number of boosting rounds and directly impacts how many trees are built in the ensemble. The best parameter set identified through hyperparameter tuning was a depth of 8, number of iterations equal to 300, and a learning rate of 0.1. The tuned model achieved a best score of -64.23, which represents an improvement in minimizing prediction error compared to other tested configurations.

The modeling section of HSMLM introduces a diverse set of base learners whose hyperparameters were selected based on the best-performing configurations obtained through GridSearchCV for each individual model. The KNN regressor predicts values by averaging the outcomes of the five closest neighbors (n_neighbors = 5). The DT algorithm is restricted to a maximum depth of ten (max_depth = 10) to avoid overfitting, while the RFR consists of 400 trees (n_estimators = 400), each trained on random subsets of the data to enhance robustness. The XGBoost model applies gradient boosting with 600 boosting rounds (n_estimators = 600), a conservative learning rate of 0.03 to ensure gradual updates, and a maximum tree depth of six (max_depth = 6), while incorporating subsampling (subsample = 0.9) and column sampling (colsample_bytree = 0.9) to improve generalization. The CatBoost is optimized with 1200 boosting iterations, a tree depth of eight, and a learning rate of 0.03, along with an l2_leaf_reg = 6.0 regularization term to control complexity and reduce overfitting.

To combine the strengths of these models, a Stacking Regressor is employed, where the predictions of the base models are passed to a meta-learner. The chosen meta-model is an ElasticNetCV, a linear regression method that applies both L1 and L2 regularization. The parameter l1_ratio = [0.1, 0.5, 0.9] specifies different balances between L1 and L2 penalties,

while alphas = np.logspace(−4, 1, 12) tests a range of penalty strengths. Cross-validation with five folds (cv = 5) ensures that the best set of hyperparameters is selected. Additionally, the passthrough = True argument ensures that the meta-model not only sees the predictions of the base models but also the original features, enriching the final prediction process.

Finally, the complete modeling workflow is wrapped in a pipeline, which integrates preprocessing and stacking into a single object. To address skewness in the target variable, the model is trained inside a Transformed Target Regressor that applies a Yeo–Johnson power transformation, making the target distribution more symmetric and thus easier for the models to learn. The fit method at the end trains the pipeline end-to-end, so that preprocessing, base model training, stacking, and target transformation are executed in a seamless and unified manner.

In the constructed models DT, XGBoost, KNN, RFR, CatBoost and HSMLM, each showed strong learning ability on the training dataset, producing low error values and high accuracy. Their performance on the training data is illustrated in Figure 8. However, the true reliability of a model is best determined by how it performs on the test dataset. Test data results act as the key benchmark, showing how effectively the models can uncover real patterns and provide accurate predictions for unseen cases. Therefore, this study places greater importance on test data performance, as it more clearly reflects the models' generalization ability and practical usefulness.

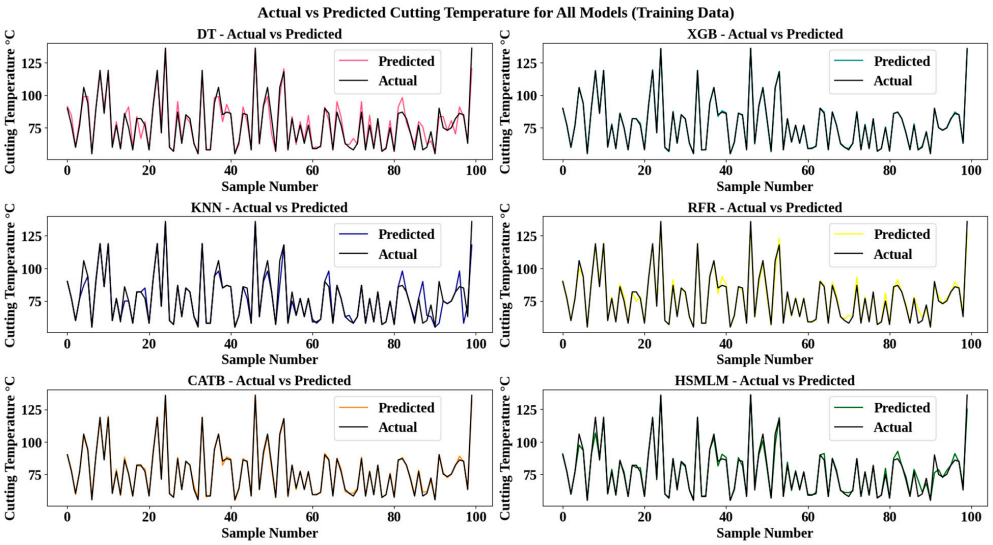

Figure 8. Predictive performances of models with training data.

The prediction performance of all ML models on the testing dataset was evaluated using multiple statistical metrics. MAE, RMSE, and MAPE are presented in Figure 9, MSE and Accuracy in Figure 10, and R^2 in Figure 11. The DT model demonstrated reliable predictive capability, achieving an MAE of 4.26, an RMSE of 6.04, an MAPE of 5.4, an MSE of 36.4, an accuracy of 94.5%, and an R^2 of 0.88. Its relatively strong performance can be attributed to its ability to capture non-linear relationships and interactions between input features efficiently, allowing it to model complex patterns in the data [61]. However, despite these strengths, the relatively simple tree structure may still limit its generalization ability in certain cases.

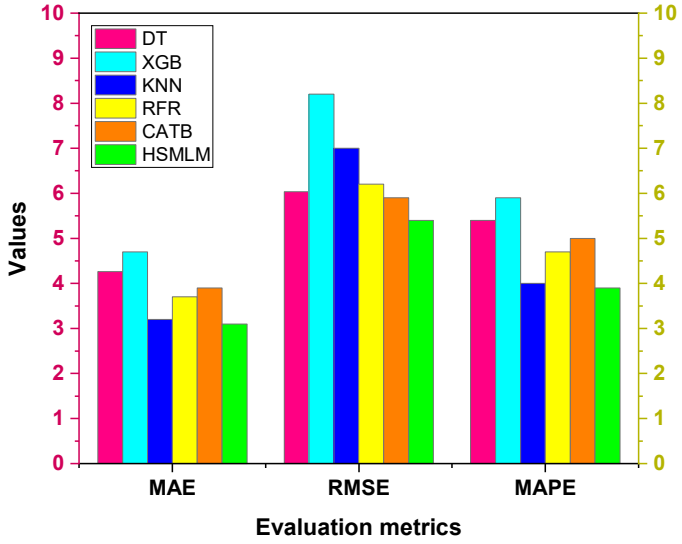

Figure 9. MAE, RMSE and MAPE of ML models.

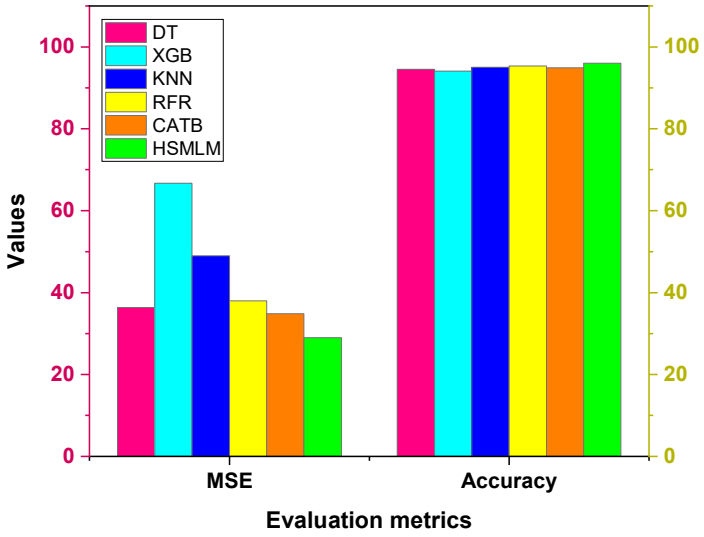

Figure 10. MSE and Accuracy of ML models.

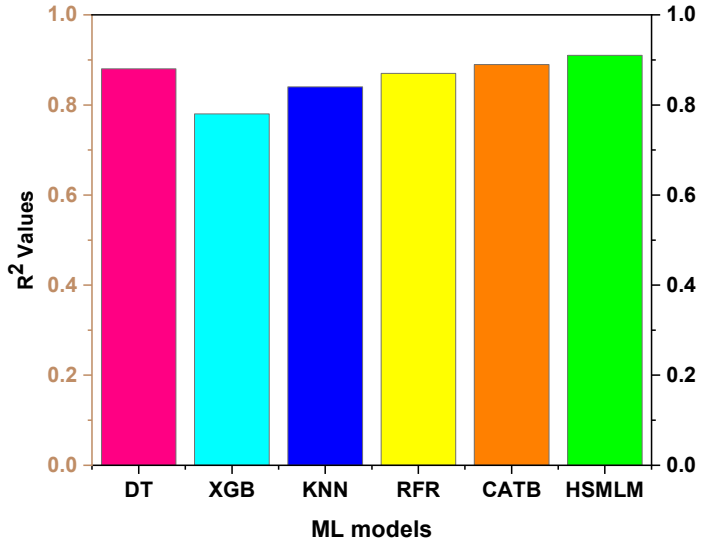

Figure 11. R^2 values of ML models.

The KNN model produced lower MAE (3.2) and MAPE (4.0) compared with DT, indicating stronger local prediction accuracy. However, its RMSE (7.0) and MSE (49) were higher, reflecting sensitivity to data variability. The accuracy (95%) was comparable to DT, while the R^2 value (0.84) was slightly lower, indicating reduced overall explanatory power. The comparatively lower performance in terms of global error metrics can be attributed to KNN's reliance on local neighbors, which makes it sensitive to fluctuations in the data and less effective at capturing overall trends across the entire dataset [62].

The RFR achieved an MAE of 3.7, an RMSE of 6.2, an MAPE of 4.7, an MSE of 38, an accuracy of 95.3%, and an R^2 of 0.87. Its high performance is due to the ensemble design, which aggregates multiple decision trees to lower variance and enhance robustness while keeping bias at a moderate level. This allows RFR to capture patterns more effectively than individual trees and generalize well across the dataset [63]. The CatBoost model outperformed RFR in terms of RMSE (5.9), MSE (34.8), and R^2 (0.89), demonstrating superior capability in modeling complex non-linear relationships and interactions within the data. However, MAE (3.9), MAPE (5.0), and accuracy (94.9%) were slightly lower, which may be due to the algorithm's sensitivity to certain localized variations in the dataset and its reliance on gradient boosting iterations [64], making it occasionally less robust to small fluctuations compared with RFR.

The XGBoost model showed the lowest performance among all models, with a MAE of 4.7, RMSE of 8.2, MAPE of 5.9, MSE of 66.7, accuracy of 94.1%, and R^2 of 0.78. The reduced performance can be attributed to its sensitivity to hyperparameter settings and higher variance, which may limit its ability to generalize effectively on the test dataset, particularly in regions with complex or sparse data patterns [65]. The proposed HSMLM achieved the best performance, with an MAE of 3.1, an RMSE of 5.4, an MAPE of 3.9, an MSE of 29, an accuracy of 96%, and an R^2 of 0.91. Its superior performance can be attributed to the ensemble stacking approach, which effectively combines the strengths of individual base models, reducing both bias and variance, and enhancing predictive robustness and generalization to unseen data. Correspondingly, Zeng et al. reported strong predictive capability using a physics-informed meta-learning model, achieving an R^2 of 0.89 and MAPE of 12.33 for surface roughness prediction in the milling of S45C steel [66]. Figure 12 presents a comparison of the predictive performance of all six models evaluated on the testing dataset.

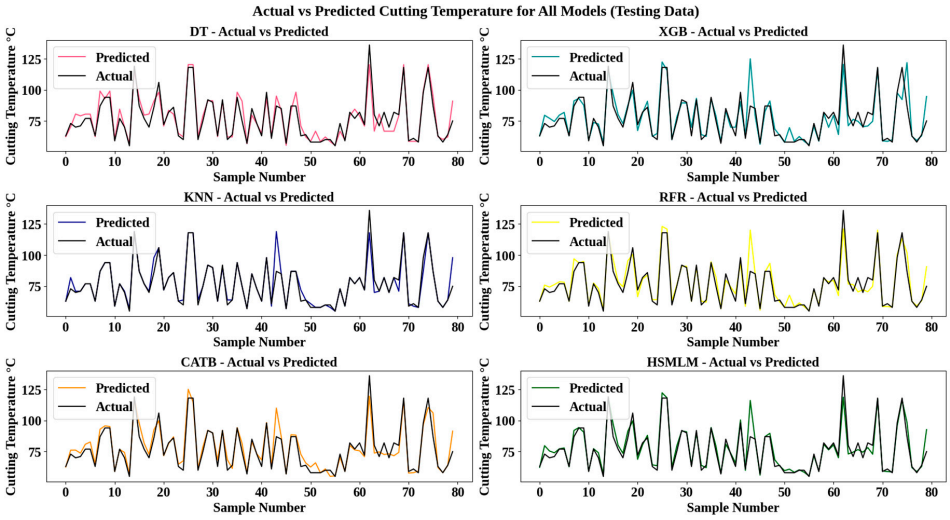

Figure 12. Predictive performances of models with testing data.

4. Conclusions

This study examines the effect of Gnps-enhanced sesame oil in an MQL system during the milling of H11 hot die steel. An HSMLM model was devised to predict cutting temperature, and its performance was compared with base learners, including DT, XGBoost, KNN, RFR, and CatBoost. The following conclusions are drawn:

- The GEMQL environment notably lowers T_c compared to CMQL, due to Gnps-enhanced lubrication that improves thermal stability, fluid retention, and protective tribo-film formation at the tool–workpiece interface.
- The lowest T_c of 58 °C in the CMQL environment was observed at a V_c of 40 m/min, a f of 0.01 mm/rev, and a P_j of 4 bar.
- The GEMQL environment achieved a maximum T_c reduction of 36.8% compared to CMQL at a V_c of 60 m/min, a f of 0.01 mm/rev, and a P_j of 2 bar. This reduction surpasses the result reported in the existing study employing Gnps-enhanced sunflower oil under MQL milling.
- Under the optimum parameters of 40 m/min V_c, 0.03 mm/rev f, and 4 bar P_j, GEMQL achieves the lowest T_c of 55 °C, significantly lower than CMQL.
- Among the six ML models, XGBoost showed the lowest performance (MAE: 4.7, RMSE: 8.2, MAPE: 5.9, MSE: 66.7, accuracy: 94.1%, R^2: 0.78) due to its sensitivity to hyperparameters and higher variance, limiting generalization on sparse test data.
- The proposed HSMLM achieved the highest performance with an MAE of 3.1, an RMSE of 5.4, an MAPE of 3.9, an MSE of 29, an accuracy of 96%, and an R^2 of 0.91, as its ensemble stacking effectively combines base models to enhance accuracy, robustness, and generalization. It also performs better than the existing physics-informed meta-learning approach reported for surface roughness prediction.

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Conflicts of Interest: The authors declare no conflict of interest.

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