

Forecasting Stock Price Accuracy for Vietnamese Pharmaceutical Firms through Ensemble Statistical, Machine Learning, and Deep Learning Models

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ABSTRACT Stock price prediction plays a crucial role in investment decision-making and portfolio management strategies. This study aims to forecast the stock prices of three major pharmaceutical companies listed on the Vietnamese stock exchange: **Hataphar (DHT)**, **Central Pharmaceutical JSC No3 (DP3)**, and **American Vietnamese Biotech Inc (AMV)**.

The research employs a ensemble of statistical and machine learning models, including Fast Fourier Transform (FFT), Linear Regression with Calendar/Fourier Terms, Deterministic Process Models, LightGBM, ARIMA, Recurrent Neural Networks (RNNs), Gated Recurrent Units (GRUs), and Long Short-Term Memory (LSTM) networks. These diverse techniques are utilized to capture the complex patterns and dynamics inherent in stock price movements.

The research evaluates model performance and identifies key factors influencing price fluctuations in Vietnam's pharmaceutical sector. Findings provide insights for investment strategies and portfolio optimization in this industry. The accurate forecasting models enable informed decision-making by investors, analysts and stakeholders.

INDEX TERMS Vietnamese stock market, Machine learning models, Statistical models, LightGBM, Recurrent neural networks, LSTM, GRU, ARIMA, Linear regression, Calendar/Fourier terms, pharma companies, AMV, DP3, DHT, FFT, Deterministic process models.

I. INTRODUCTION

The Covid-19 pandemic has severely disrupted global economies, with the pharmaceutical industry at the forefront combating this crisis. As demand for vaccines, treatments, and medical supplies spiked, pharmaceutical firms faced immense pressures, leading to significant volatility in their stock prices. In Vietnam, companies like **Hataphar**, **Central Pharmaceutical No3**, and **American Vietnamese Biotech** played vital roles supplying Covid-19 countermeasures amidst shifting market dynamics.

Traditional forecasting models may not fully capture the unprecedented pandemic's complexities impacting stock movements. This necessitates exploring advanced machine learning and deep learning techniques to effectively model intricate patterns and non-linearities. This study aims to leverage an ensemble of statistical methods (FFT, Linear Regression with Calendar/Fourier terms, ARIMA) and machine learning models (LightGBM, RNNs, GRUs, LSTMs) for forecasting stock prices of the three major Vietnamese

pharmaceutical firms.

By combining diverse approaches, the research seeks insights into factors influencing price dynamics during Covid-19. Findings can contribute to financial forecasting knowledge, guiding investment strategies, portfolio optimization, and risk management for stakeholders in Vietnam's pharmaceutical sector. Reliable forecasts can support the industry's resilience and growth amidst future crises.

II. RELATED WORKS

Xiao and Wang (2017) [1] applied ARIMA, LSTM and a hybrid ARIMA-LSTM model to forecast stock prices of two companies listed on Chinese stock exchanges. Using data from 2009-2016, they found the hybrid model outperformed individual ARIMA and LSTM across evaluation metrics like RMSE and MAPE, improving accuracy by 10-15%.

Nelson et al. (2017) [2] compared statistical (ARIMA, Exponential Smoothing) and machine learning models (ANN, LSTM, GBR) for forecasting stock volatility of the S&P

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500 index. Using 60 years of daily data, they showed LSTM networks were most effective in capturing nonlinear patterns and long-range temporal dependencies.

Bao et al. (2017) [3] proposed a novel wavelet transformer combining wavelet analysis and self-attention for stock movement prediction on high-frequency intraday data of MSCI USA Index constituents in 2017. Evaluated against statistical baselines, ARIMA and Exponential Smoothing, the wavelet transformer demonstrated superior directional prediction accuracy.

Minh et al. (2018) [4] explored CNN, LSTM and hybrid CNN-LSTM architectures for forecasting stock prices of major companies listed on the Vietnamese stock exchange like VNM, VCB and BVH, using data from 2012-2017. Their CNN-LSTM model incorporating technical indicators outperformed individual CNN and LSTM networks.

Chen et al. (2020) [5] developed an adaptive, dynamic method based on time series clustering for stock forecasting. Applied to constituent stocks of S&P 500 and CSI 300 indexes from 2013-2018, their approach using K-means clustering and Exponential Smoothing outperformed classical statistical models like ARIMA.

Mamon et al. (2021) [6] conducted a comparative analysis assessing the performance of ARIMA, VAR, CNN, LSTM and hybrid CNN-LSTM models on predicting Philippine stock index PSEi returns during the Covid-19 pandemic from 2020-2021 data. The CNN-LSTM exhibited the lowest error rates across metrics.

Jolly Masih [7] conducted a study that implemented algorithms such as SVM and LSTM on stock market data to observe whether major IT companies experienced growth or decline during the COVID-19 pandemic. The authors also utilized the ARIMA forecasting method to predict the stock prices of the mentioned four companies: Google, Microsoft, Apple, and Amazon. Regarding the current prediction of stock prices, the authors achieved an accuracy of 75.95% using the SVM algorithm and the best-fit forecast model of ARIMA. This decrease in accuracy was anticipated, as described earlier, due to increased volatility in the stock market resulting from the COVID-19 outbreak.

Ke et al. (2017) [20] introduced LightGBM, a highly efficient gradient boosting decision tree algorithm for classification and regression tasks. This method has gained significant attention for its superior performance, efficient handling of high-dimensional data, and effective parallelization. Several studies have successfully applied LightGBM to stock price prediction problems, leveraging its ability to capture nonlinear relationships and feature interactions.

III. MATERIALS

A. DATASET

We have compiled three datasets in the pharmaceutical field from three companies **Hataphar** (DHT), Central Pharmaceutical JSC No3 (DP3), and American Vietnamese Biotech Inc (AMV). The collected data is the historical stock price of the three companies from 1-1-2019 to 17/4/2024. The data was collected from Investing.com website and includes columns:

TABLE 1. Data's attributes meaning

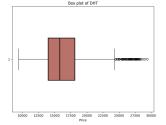
	ъ		
Attributes	Description		
Date	The date when the stock price data		
	was recorded.		
Price	Closing price of the stock for each		
	date.		
Open	Opening price of the stock for each		
	date.		
High	Highest price of the stock during		
	the trading session.		
Low	Lowest price of the stock during the		
	trading session.		
Vol.	Trading volume of the stock on		
	each date.		
Change %	Percentage change in the stock		
	price compared to the previous trad-		
	ing day's closing price.		

The objective is to predict the price (close price), so only descriptive statistical data relating to the "Price" column will be processed.

B. DESCRIPTIVE STATISTICS

TABLE 2. DHT, DP3, AMV's Descriptive Statistics

	DHT	DP3	AMV
Count	1320	1146	1322
Mean	16,583	86,519	11,167
Std	4,149	21,584	5,811
Min	9,384	57,600	2,700
25%	13,961	68,000	4,900
50%	15,778	81,000	11,359
75%	18,133	103,000	15,123
Max	29,400	157,500	24,589



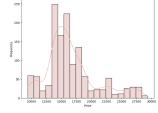
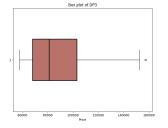


FIGURE 1. DHT stock price's boxplot

FIGURE 2. DHT stock price's histogram



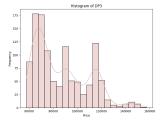
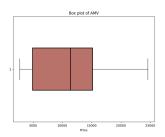


FIGURE 3. DP3 stock price's boxplot

FIGURE 4. DP3 stock price's histogram





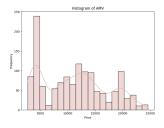


FIGURE 5. AMV stock price's boxplot

FIGURE 6. AMV stock price's histogram

IV. METHODOLOGY

A. LINEAR REGRESSION

Regression analysis is a tool for building mathematical and statistical models that characterize relationships between a dependent variable and one or more independent, or explanatory, variables, all of which are numerical. This statistical technique is used to find an equation that best predicts the y variable as a linear function of the x variables. A multiple linear regression model has the form:

$$Y = \beta_0 + \beta_1 X_1 + \beta_2 X_2 + \dots + \beta_k X_k + \varepsilon$$

Where:

- Y is the dependent variable (Target Variable).
- X_1, X_2, \ldots, X_k are the independent (explanatory) variables
 - β_0 is the intercept term.
- $\beta_1, ..., \beta_k$ are the regression coefficients for the independent variables.
 - ε is the error term.

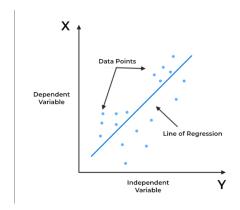


FIGURE 7. LN model's

B. FAST FOURIER TRANSFORM

The Fast Fourier Transform (FFT) is a method that allows computing the Discrete Fourier Transform (DFT) in $\mathcal{O}(n\log n)$ time. The basic idea of the FFT is to apply the divide-and-conquer approach. We divide the coefficient vector of the polynomial into two vectors, recursively compute the DFT for each of them, and combine the results to compute the DFT of the complete polynomial.

Let there be a polynomial A(x) with degree n-1, where n is a power of 2, and n>1:

$$|A(x)| = a_0 x^0 + a_1 x^1 + \dots + a_{n-1} x^{n-1}$$

Where:

- |A(x)| is the absolute value of the polynomial A(x).
- $a_0, a_1, \ldots, a_{n-1}$ are the coefficients of the polynomial.
- x^0, x^1, \dots, x^{n-1} are the corresponding powers of the variable x for each coefficient.
- n-1 is the highest degree of the polynomial.
- The · · · (read as "and so on") represents the remaining terms in the sequence.

We divide it into two smaller polynomials, one containing only the coefficients of the even positions, and the other containing the coefficients of the odd positions:

$$A_0(x) = a_0 x^0 + a_2 x^1 + \dots + a_{n-2} x^{\frac{n}{2} - 1}$$

$$A_1(x) = a_1 x^0 + a_3 x^1 + \dots + a_{n-1} x^{\frac{n}{2} - 1}$$

In the FFT computation, we apply the divide-and-conquer approach recursively to the polynomials $A_0(x)$ and $A_1(x)$, and combine the results to obtain the DFT of the original polynomial A(x). Dividing the polynomial into smaller polynomials and computing recursively significantly reduces the computational time compared to the traditional DFT computation method.

C. ARIMA

The ARIMA model (an acronym for Auto-Regressive Integrated Moving Average), essentially creates a linear equation which describes and forecasts your time series data. This equation is generated through three separate parts which can be described as:

- AR auto-regression: equation terms created based on past data points.
- I integration or differencing: accounting for overall "trend" in the data.
- MA moving average: equation terms of error or noise based on past data points.

An ARIMA model is characterized by 3 terms: p, d, q where,

- p is the order of the AR term.
- q is the order of the MA term.
- d is the number of differencing required to make the time series stationary.

A pure Auto Regressive (AR only) model is one where Yt depends only on its own lags. That is, Yt is a function of the 'lags of Yt'.

$$Y_t = a + B_1 Y_{t-1} + B_2 Y_{t-2} + \ldots + B_p Y_{t-p} + \varepsilon_t$$

Where

- Y_t : The value of the time series at time t.
- a: A constant term.
- B_1, B_2, \dots, B_p : Autoregressive coefficients representing the relationship between the current observation and its past values at lag intervals.



- p: The order of the autoregressive component.
- ε_t : A white noise error term at time t.

Likewise a pure Moving Average (MA only) model is one where Yt depends only on the lagged forecast errors.

$$Y_t = a + \phi_1 \varepsilon_{t-1} + \phi_2 \varepsilon_{t-2} + \ldots + \phi_q \varepsilon_{t-q} + \varepsilon_t$$

Where

• q: The order of the moving average component.

An ARIMA model is one where the time series was differenced at least once to make it stationary and you combine the AR and the MA terms. So the equation becomes:

$$Y_t = a + \sum_{i=1}^{p} \phi_i Y_{t-i} + \varepsilon_t + \sum_{j=1}^{q} B_j \varepsilon_{t-j}$$

To find the orders of AR, I, and MA components in an ARIMA model, analysts use ACF and PACF plots. A decreasing ACF suggests a MA component, with the lag where it drops below a threshold indicating its order. Significant spikes in PACF imply an AR component, with the lag where PACF cuts off after a spike revealing its order. By interpreting these plots, analysts select suitable orders for ARIMA modeling.

D. LSTM

Long Short-Term Memory (LSTM) network is a type of recurrent neural network (RNN) capable of learning long-term dependencies. The LSTM model consists of several components, including gates that control the flow of information through the network. The LSTM model can be described as follows:

- Let \mathbf{x}_t be the input vector at time step t.
- Let \mathbf{h}_t be the hidden state vector at time step t.
- Let c_t be the cell state vector at time step t.
- Let W, U, and b be the weight matrices and bias

The LSTM model is defined by the following equations:

1. Forget Gate:

$$\mathbf{f}_t = \sigma(\mathbf{W}_f \mathbf{x}_t + \mathbf{U}_f \mathbf{h}_{t-1} + \mathbf{b}_f)$$

2. Input Gate:

$$\mathbf{i}_t = \sigma(\mathbf{W}_i \mathbf{x}_t + \mathbf{U}_i \mathbf{h}_{t-1} + \mathbf{b}_i)$$

3. Candidate Memory Cell:

$$\mathbf{\tilde{c}}_t = \tanh(\mathbf{W}_c \mathbf{x}_t + \mathbf{U}_c \mathbf{h}_{t-1} + \mathbf{b}_c)$$

4. Cell State:

$$\mathbf{c}_t = \mathbf{f}_t \odot \mathbf{c}_{t-1} + \mathbf{i}_t \odot \tilde{\mathbf{c}}_t$$

5. Output Gate:

$$\mathbf{o}_t = \sigma(\mathbf{W}_o \mathbf{x}_t + \mathbf{U}_o \mathbf{h}_{t-1} + \mathbf{b}_o)$$

6. Hidden State:

$$\mathbf{h}_t = \mathbf{o}_t \odot \tanh(\mathbf{c}_t)$$

Where:

- σ denotes the sigmoid activation function.
- tanh denotes the hyperbolic tangent activation function.
- ① denotes element-wise multiplication.
- $\mathbf{W}_f, \mathbf{W}_i, \mathbf{W}_c, \mathbf{W}_o$ are the input weight matrices.
- $\mathbf{U}_f, \mathbf{U}_i, \mathbf{U}_c, \mathbf{U}_o$ are the recurrent weight matrices.
- $\mathbf{b}_f, \mathbf{b}_i, \mathbf{b}_c, \mathbf{b}_o$ are the bias vectors.
- \mathbf{f}_t is the forget gate vector.
- \mathbf{i}_t is the input gate vector.
- $\tilde{\mathbf{c}}_t$ is the candidate memory cell vector.
- \mathbf{c}_t is the cell state vector.
- **o**_t is the output gate vector.
- \mathbf{h}_t is the hidden state vector.

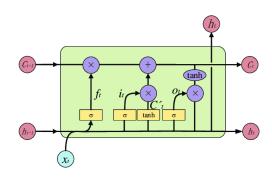


FIGURE 8. LSTM model's

E. LINEAR REGRESSION COMBINED WITH CALENDARFOURIER AND DETERMINISTICPROCESS

Linear Regression combined with CalendarFourier and DeterministicProcess is an extended linear regression method used for modeling and forecasting time series with recurring cycles and growth trends. This method combines the following elements:

Linear Regression:

- Linear Regression is a common statistical modeling technique used to explore the linear relationship between an output variable (dependent variable) and one or more input variables (independent variables).
- Linear Regression is often used as a foundation for other time series forecasting techniques due to its ability to effectively model linear relationships.

CalendarFourier:

- CalendarFourier is a time feature used to model recurring patterns in time series data.
- It transforms time into Fourier functions (sine and cosine), allowing the model to learn cyclical patterns such as daily, weekly, monthly, or yearly cycles.
- When combined with Linear Regression, Calendar-Fourier enhances the ability to forecast recurring cycles in time series data.

DeterministicProcess:

• DeterministicProcess is a time feature used to model trends or growth patterns in time series data.



- It transforms time into a growth function (e.g., straight line, curve, polynomial, etc.), allowing the model to learn the long-term growth or decline trends of the data.
- When combined with Linear Regression, DeterministicProcess improves the ability to forecast long-term trends in time series data.

Suppose we have a time series y and input variables X. The Linear Regression model combined with CalendarFourier and DeterministicProcess takes the form:

$$y = \beta_0 + \beta_1 X_1 + \beta_2 X_2 + \dots + \beta_n X_n + f(t) + g(t) + \epsilon$$

Where:

- y is the output variable (the time series to be predicted).
- β_0 is the constant term.
- $\beta_1, \beta_2, ..., \beta_n$ are the regression coefficients for the input variables $X_1, X_2, ..., X_n$.
- f(t) is the CalendarFourier component, representing the recurring cycles in the time series, with t being the time.
- g(t) is the DeterministicProcess component, representing the growth trend in the time series, with t being the time.
- ϵ is the random error term.

The CalendarFourier component f(t) can be represented as a sum of sine and cosine functions with different cycle frequencies:

$$f(t) = \alpha_1 \sin(\omega t) + \beta_1 \cos(\omega t) + \alpha_2 \sin(2\omega t) + \beta_2 \cos(2\omega t) + \dots + \beta_3 \sin(2\omega t) + \beta_3 \cos(2\omega t) + \dots + \beta_3 \sin(2\omega t) + \beta_3 \cos(2\omega t) + \dots + \beta_3 \cos(2\omega t) + \beta_3 \cos(2\omega t) + \dots + \beta_3 \cos(2\omega t)$$

$$\alpha_k \sin(k\omega t) + \beta_k \cos(k\omega t)$$

Where:

- ω is the cycle frequency (e.g., $2\pi/365$ for an annual cycle)
- k is the number of Fourier terms used to model the cycles.
- α₁, β₁, α₂, β₂, ..., α_k, β_k are the coefficients to be estimated from the data.

The DeterministicProcess component g(t) can be represented by a growth function, such as a polynomial, exponential, or logarithmic function:

$$g(t)=\gamma_1 t + \gamma_2 t^2 + \gamma_3 t^3 + \dots$$
 (polynomial)
$$g(t)=\gamma_1 + \gamma_2 \log(t) \quad \text{(logarithmic)}$$

$$g(t)=\gamma_1 \exp(\gamma_2 t) \quad \text{(exponential)}$$

Where $\gamma_1, \gamma_2, \gamma_3, ...$ are the coefficients to be estimated from the data.

The estimation of the coefficients $\beta_0, \beta_1, \beta_2, ..., \beta_n, \alpha_1, \beta_1, \alpha_2$, can be performed using the Ordinary Least Squares (OLS) method or other optimization techniques.

By combining these components, the Linear Regression model with CalendarFourier and DeterministicProcess can effectively model and forecast both recurring cycles and growth trends in the time series, thereby achieving higher forecasting accuracy compared to simpler models.

F. GATED RECURRENT UNIT

The Gated Recurrent Unit (GRU) is a type of recurrent neural network (RNN) architecture similar to Long Short-Term Memory (LSTM) but with a simpler design and fewer parameters. GRU employs two gates: the reset gate and the update gate. The reset gate controls the forgetting of the previous hidden state, while the update gate manages the incorporation of a candidate activation vector into the new hidden state.

The GRU processes sequential data one element at a time, updating its hidden state based on the current input and the previous hidden state. At each time step, it computes a "candidate activation vector" that combines information from the input and previous hidden state. This vector is updated using the reset gate, which controls how much of the previous hidden state to forget, and the update gate, which determines how much of the candidate activation vector to incorporate into the new hidden state.

Here's the math behind the GRU architecture:

1. The reset gate r and update gate z are computed using the current input x and the previous hidden state h_{t-1} :

$$r_t = \operatorname{sigmoid}(\mathbf{W}_r \cdot [h_{t-1}, x_t])$$

$$z_t = \operatorname{sigmoid}(\mathbf{W}_z \cdot [h_{t-1}, x_t])$$

where \mathbf{W}_r and \mathbf{W}_z are weight matrices that are learned during training.

2. The candidate activation vector \tilde{h}_t is computed using the current input x and a modified version of the previous hidden state that is "reset" by the reset gate:

$$\tilde{h}_t = \frac{e^{r_t * h_{t-1}} + e^{-r_t * h_{t-1}}}{e^{r_t * h_{t-1}} - e^{-r_t * h_{t-1}}} \cdot x_t$$

where \mathbf{W}_h is another weight matrix.

3. The new hidden state h_t is computed by combining the candidate activation vector with the previous hidden state, weighted by the update gate:

$$h_t = (1 - z_t) * h_{t-1} + z_t * \tilde{h}_t$$

Overall, the reset gate controls the retention of the previous hidden state, and the update gate determines how much of the candidate, activation, vector updates the new hidden state. This compact architecture selectively updates its hidden state without needing a separate memory cell like LSTM.



G. LIGHTGBM

LightGBM, short for Light Gradient Boosting Machine, is a gradient boosting framework developed by Microsoft. It's renowned for its efficiency and speed. The fundamental representation of how LightGBM combines predictions from different trees to make a final prediction can be represented by the following formula:

$$Y = \text{Base_tree}(X) - \text{lr} \times \text{Tree}_1(X) - \text{lr} \times \text{Tree}_2(X) - \text{lr} \times \text{Tree}_3(X)$$

Here, Y is the final prediction, Base tree(X) is the prediction of the base model, Tree1(X), Tree2(X), and Tree3(X) are the predictions of subsequent models, and lr is the learning rate

LightGBM introduces two key techniques: Gradient-based One-Side Sampling (GOSS) and Exclusive Feature Bundling (EFB).

- GOSS focuses on sampling instances based on their gradients, prioritizing those with larger gradients for training. It also randomly drops instances with small gradients, thereby reducing the computational burden.
- EFB optimizes memory usage by bundling exclusive features together, allowing for more efficient processing and faster training without compromising the performance significantly.

LightGBM employs histogram-based algorithms to find the best split points during tree construction1. Instead of exhaustively searching through all possible split points, Light-GBM discretizes continuous features into discrete bins and constructs histograms to approximate the optimal split points. This approach significantly reduces computational complexity and memory usage, especially for datasets with a large number of unique feature values.

LightGBM leaf-wise

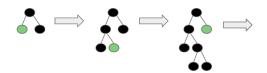


FIGURE 9. LightGBM model's

H. RECURRENT NEURAL NETWORK

Recurrent Neural Network (RNN) is a neural network architecture that enables processing sequential data in O(n) time complexity. The core idea of RNNs is to apply the divide-and-conquer approach. We divide the input sequence into time steps, recursively compute for each step, and combine the results to process the complete input sequence. Let X be an input sequence of length n:

$$|X| = x_1, x_2, ..., x_n \tag{1}$$

Where:

• |X| is the length of the input sequence X.

- $x_1, x_2, ..., x_n$ are the elements of the sequence.
- ... (read as "and so on") represents the remaining elements in the sequence.

We divide the input sequence into time steps, processing one element at each step:

$$h_1 = f(x_1, h_0) h_2 = f(x_2, h_1) \dots h_n = f(x_n, h_{n-1})$$
(2)

In which:

- h_t is the hidden state at time step t.
- f is a computation function, typically a non-linear transformation (e.g., tanh, ReLU) of the linear combination of the input x_t and the previous hidden state h_{t-1} .
- h_0 is the initial hidden state, usually initialized as zeros.

In the RNN computation, we apply the divide-and-conquer approach recursively at each time step, computing the hidden state h_t based on the input x_t and the previous hidden state h_{t-1} . The final result is the hidden state h_n , which encapsulates the contextual information from the entire input sequence.

The output y_t at each time step t is computed based on the current hidden state h_t and input x_t , typically through an output layer:

$$y_t = g(h_t, x_t) \tag{3}$$

Where g is a function, often a linear transformation or a neural network.

Recurrent structure allows RNNs to capture long-range dependencies within the sequential data, making them suitable for various tasks such as language modeling, machine translation, speech recognition, and time series prediction.

V. RESULT

A. EVALUATION METHODS

Mean Percentage Absolute Error (MAPE): is the average percentage error in a set of predicted values.

$$MAPE = \frac{100\%}{n} \sum_{i=1}^{n} |y_i - \hat{y}_i| = 1$$

Root Mean Squared Error (RMSE): is the square root of average value of squared error in a set of predicted values.

$$RMSE = \sqrt{\sum_{i=1}^{n} \frac{(\hat{y_i} - y_i)^2}{n}}$$

Mean Absolute Error (MSLE):is the relative difference between the log-transformed actual and predicted values.

$$MSLE = \frac{1}{n} \sum_{i=1}^{n} (log(1 + \hat{y}_i) - log(log(1 + y_i))^2)$$

Where:

- \bullet n is the number of observations in the dataset.
- y_i is the true value.
- $\hat{y_i}$ is the predicted value.



B. AMV DATASET

AMV Dataset's Evaluation				
Model	Training:Testing	RMSE	MAPE (%)	MSLE
LN	7:3	7781.26	49.62	0.19
	8:2	9119.01	67.34	0.27
	9:1	6907.31	51.82	0.17
RNN	7:3	1276.19	0.3207	0.0801
KININ	8:2	135.78	0.0264	0.0011
	9:1	264.547	0.0665	0.0054
GRU	7:3			
J GKU	8:2			
	9:1			
ADDAA	7:3	558.44	11.50%	0.02
ARIMA	8:2	801.64	18.26%	0.04
	9:1	290.25	5.51%	0.00
FFT	7:3	0.0829	4151	0.0054
LL I	8:2	0.05950	0.9657	0.003
	9:1	0.0259	0.5279	0.0006
LN-CF-DP	7:3	6841.46	165.18	0.96
LN-CF-DP	8:2	2031.85	44.21	0.15
	9:1	1441.58	33.21	0.10
LSTM	7:3	1861.67	1.78%	0.00068
LSTW	8:2	522.86	2.41%	0.00135
	9:1	298.001	1.74%	0.00046
LightGRM	7:3	0.11	215.22	0.01
LightGBM	8:2	0.05	128.70	0.002
	9:1	0.05	188.43	0.002

TABLE 3. AMV Dataset's Evaluation

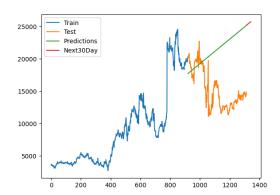


FIGURE 10. Linear model's result with 7:3 splitting proportion

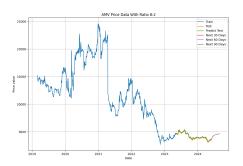


FIGURE 11. RNN model's result with 8:2 splitting proportion

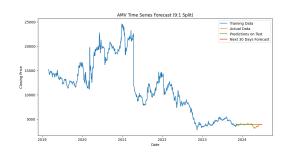


FIGURE 12. Arima model's result with 8:2 splitting proportion

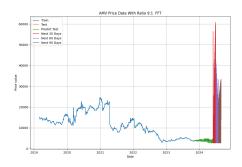


FIGURE 13. FFT model's result with 9:1 splitting proportion

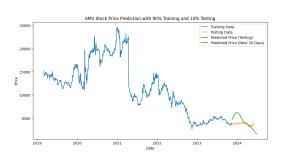


FIGURE 14. LN-CF-DP model's result with 9:1 splitting proportion

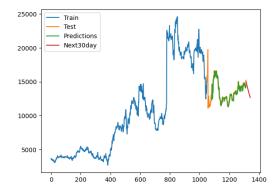


FIGURE 15. LSTM model's result with 8:2 splitting proportion



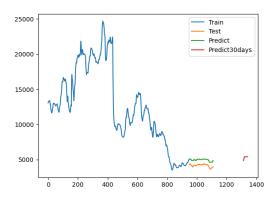


FIGURE 16. LightGBM model's result with 8:2 splitting proportion

FIGURE 18. RNN model's result with 9:1 splitting proportion

C. DP3 DATASET

	DP3 Dataset's Evaluation			
Model	Training:Testing	RMSE	MAPE (%)	MSLE
LN	7:3	7781.26	49.62	0.19
	8:2	9119.01	67.34	0.27
	9:1	6907.31	51.82	0.17
RNN	7:3	1675.09	0.0412	0.0037
KININ	8:2	2503.94	0.0700	0.0081
	9:1	1499.11	0.0275	0.0021
GRU	7:3			
GKU	8:2			
	9:1			
ADDAA	7:3	27750.14	39.01%	0.36
ARIMA	8:2	7936.36	10.027%	0.01
	9:1	6394.82	7.79%	0.01
FFT	7:3	0.2076	61.27%	0.0232
	8:2	0.1987	59.18%	0.0224
	9:1	0.1352	37.86%	0.0103
I M CE DD	7:3	54681.37	69.13	0.31
LN-CF-DP	8:2	48774.18	67.56	0.28
	9:1	45471.00	66.45	0.26
LSTM	7:3	1861.67	1.78%	0.00068
LSIM	8:2	522.86	2.41%	0.00135
	9:1	298.001	1.74%	0.00046
LightCDM	7:3	0.11	215.22	0.01
LightGBM	8:2	0.05	128.70	0.002
	9:1	0.05	188.43	0.002

TABLE 4. DP3 Dataset's Evaluation

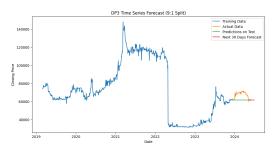


FIGURE 19. ARIMA model's result with 8:2 splitting proportion

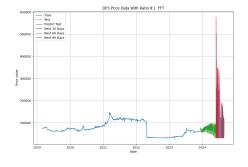


FIGURE 20. FFT model's result with 9:1 splitting proportion

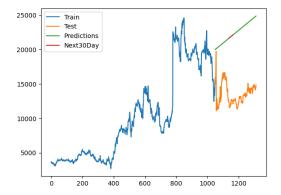


FIGURE 17. Linear model's result with 8:2 splitting proportion



FIGURE 21. LN-CF-DP model's result with 9:1 splitting proportion



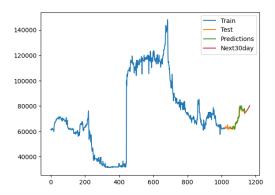


FIGURE 22. LSTM model's result with 9:1 splitting proportion

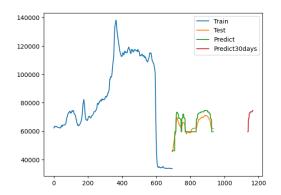


FIGURE 23. LightGBM model's result with 7:3 splitting proportion

D. DHT DATASET

DHT Dataset's Evaluation				
Model	Training:Testing	RMSE	MAPE (%)	MSLE
LN	7:3	7781.26	49.62	0.19
	8:2	9119.01	67.34	0.27
	9:1	6907.31	51.82	0.17
RNN	7:3	3065.26	0.0401	0.0026
KININ	8:2	2855.37	0.0276	0.0018
	9:1	2327.70	0.0319	0.0011
GRU	7:3			
GKU	8:2			
	9:1			
ARIMA	7:3	9640.50	29.78%	0.24
AKIMA	8:2	9305.45	31.50%	0.18
	9:1	5570.55	12.58%	0.04
FFT	7:3	0.5714	63.88%	0.0871
LL1	8:2	0.5188	57.93%	0.0862
	9:1	0.3351	34.96%	0.3354
LN-CF-DP	7:3	5726.49	24.00	0.08
LN-CF-DP	8:2	8043.72	29.24	0.16
	9:1	6725.16	24.92	0.10
LSTM	7:3	1861.67	1.78%	0.00068
LSTM	8:2	522.86	2.41%	0.00135
	9:1	298.001	1.74%	0.00046
LightCDM	7:3	0.11	215.22	0.01
LightGBM	8:2	0.05	128.70	0.002
	9:1	0.05	188.43	0.002

TABLE 5. DHT Dataset's Evaluation

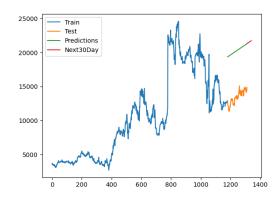


FIGURE 24. Linear model's result with 9:1 splitting proportion

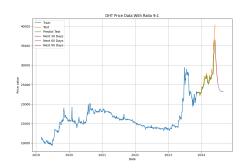


FIGURE 25. RNN model's result with 9:1 splitting proportion

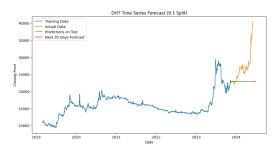


FIGURE 26. Arima model's result with 8:2 splitting proportion

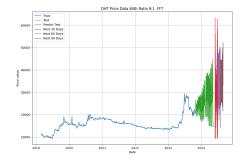


FIGURE 27. FFT model's result with 9:1 splitting proportion





FIGURE 28. LN-CF-DP model's result with 7:3 splitting proportion

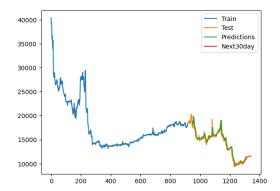


FIGURE 29. LSTM model's result with 7:3 splitting proportion

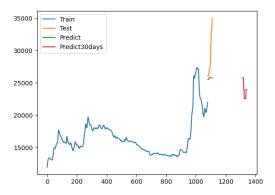


FIGURE 30. LightGBM model's result with 8:2 splitting proportion

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VI. CONCLUSION

A. SUMMARY

This research explored various methods for predicting stock prices, ranging from traditional statistical models to advanced machine learning algorithms. The models tested include Linear Regression (LR), ARIMA, GRU, RNN, LSTM, FFT, LightGBM, and LN-CF-DP. Among these, LSTM and GRU stand out as the most effective models for stock price prediction.

Each model provides unique insights: deep learning models effectively capture non-linear patterns, traditional methods offer interpretable results, while LightGBM efficiently handles multiple data features.

Although no method can perfectly predict stock prices due to market complexities, this study provides a comprehensive toolkit for investors in the pharmaceutical sector.

B. FUTURE CONSIDERATIONS

To advance this research in the future, focus should be placed on:

- Optimizing existing models: Improve the accuracy of LightGBM, LSTM, and GRU through hyperparameter tuning and advanced data processing techniques.
- Incorporating sentiment analysis: Integrate data from pharmaceutical industry news and social media to capture the impact of market sentiment.
- Expanding data scope: Add macroeconomic indicators and pharmaceutical industry-specific data to enhance predictive capabilities.
- 4) Applying ensemble learning techniques: Explore methods such as Bagging or Boosting to combine the strengths of multiple models.
- 5) Researching new models: Monitor and experiment with the latest forecasting algorithms in machine learning and quantitative finance.

By continuously exploring and incorporating new features, data sources, and modeling techniques, we can strive to continuously optimize forecasting models and enhance the ability to predict stock prices with greater accuracy and reliability.

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