

Using Forecasting Models in Cryptocurrency Price Prediction

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ABSTRACT Predicting cryptocurrency price has become a central focus for investors and traders in the rapidly evolving landscape of digital assets. Cryptocurrencies like Bitcoin, Ethereum, and other altcoins have captured global attention, with their prices subject to fluctuations influenced by various factors, from market sentiment to technological developments. With the unprecedented volatility of these markets, the need for accurate forecasting models has never been more crucial. In this research, we conduct an in-depth analysis using various models including Linear regression (LR), Autoregressive Integrated Moving Average (ARIMA), Recurrent Neural Network (RNN), Gated Recurrent Unit (GRU), Long Short-Term Memory (LSTM), Seasonal Autoregressive Integrated Moving Average with eXogenous factors (SARIMAX), Neural Basis Expansion Analysis Time Series (NBEATS), Random Forest (RF) to predict cryptocurrency price and determine the effectiveness of each model.

INDEX TERMS Cryptocurrency, price prediction, Bitcoin, Ethereum, Linear regression, ARIMA, RNN, GRU, LSTM, SARIMAX, NBEATS, RF.

I. INTRODUCTION

Cryptocurrency, a digital form of currency, has transformed the traditional concept of money. Unlike cash, cryptocurrencies exist purely in electronic form, utilizing encryption techniques for secure transactions. Bitcoin, the pioneering cryptocurrency introduced in 2009 by Satoshi Nakamoto, paved the way for a myriad of digital currencies that followed. The decentralized nature of cryptocurrencies means they operate without a central authority, using blockchain technology to record and verify transactions across a network of computers. This decentralization provides a level of transparency and security previously unseen in financial systems. As interest and adoption continue to grow, cryptocurrencies are reshaping the landscape of finance, offering individuals around the globe new avenues for transactions, investments, and financial freedom.

By using forecasting models and machine learning algorithms to analyze historical data of Bitcoin, Ethereum, and BNB, the research objective is to provide valuable insights for investors and traders to forecast the cryptocurrency price movement. To evaluate the performance of each predictive model, we use three metrics Mean Absolute Error (MAE), Mean Absolute Percentage Error (MAPE), and Root Mean Squared Error (RMSE).

II. RELATED WORKS

Mahir Iqbal and his team [1] applied a machine learning based time series analysis in predicting the price of Bitcoin. They used ARIMA, FBProphet, and XG Boosting for time series analysis and evaluated these models by using three parameters RMSE, MAE, and R Squared (R²). After conducting the analysis, they concluded that ARIMA is the best of three with RMSE score and MAE score are 322.4 and 227.3 respectively.

Harsha Nanda Gudavalli and Khetan Venkata Ratnam Kancherla [2] identified four machine learning algorithms RF, Gradient Boosting (GB), LSTM, and GRU to make long-term predictions of cryptocurrency prices. Overall, LSTM is the most accurate model based on RMSE score of 0.01083, MSE score of 0.00011, and R Squared score of 0.80618.

Junyi Zhu [3] applied LR for Bitcoin price prediction with the accuracy rate for training RSS and testing RSS are 94.6% and 96.99% respectively.

Namrata Hemraj Gawali [4] developed predictive models to predict the price of Bitcoin using LSTM, Seasonal Autoregressive Integrated Moving Average (SARIMA), SARIMAX, and RNN. In conclusion, LSTM outperformed all other models with R Squared score of 0.9519, MAE score of 2476.68, MSE score of 10958440.29 and RMSE score of 3310.35.

Alikhan Bulatov [5] used N-BEATS trained on Bitcoin daily,

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hourly, and up-to-the-minute data in comparison with LSTM and ARIMA. In this analysis, it is concluded that N-BEATS is the best model for forecasting the Bitcoin price with the RMSE score in daily, hourly, and minute data are 308.859, 59.303, and 13.678 respectively.

III. MATERIALS

A. DATASET

This research focuses on analyzing cryptocurrency trends through three data sets of three cryptocurrencies:

- BTC/USD
- ETH/USD
- BNB/USD

Those data sets are extracted from the website investing.com with the selected period from March 1, 2019, to March 1, 2024. Each data set contains 1828 data rows and has 6 attributes:

- Price: The last transacted price in a day.
- Open: The opening price in a day.
- High: The highest price in a day.
- Low: The lowest price in a day.
- Vol: The total amount traded in a day.
- Change %: The percentage change in the closing price compared to the previous day's closing price.

B. DESCRIPTIVE STATISTICS

TABLE 1. BTC, ETH, BNB's Descriptive Statistics

| | BTC | ETH | BNB |
|-------|------------|-----------|----------|
| Count | 1828 | 1828 | 1828 |
| Mean | 25,983.05 | 1,488.75 | 212.44 |
| Std | 16,801.607 | 1,160.628 | 171.7707 |
| Min | 3,715 | 107.58 | 9.28 |
| 25% | 10,255.075 | 256.75 | 25.3325 |
| 50% | 23,637.8 | 1,573.75 | 242.4 |
| 75% | 38,686.75 | 2,152.043 | 318.675 |
| Max | 67,520 | 4,808.09 | 676.12 |

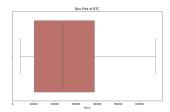
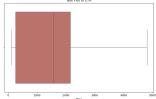


FIGURE 1. Bitcoin price's boxplot

FIGURE 2. Bitcoin price's histogram



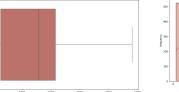


FIGURE 3. Ethereum price's boxplot

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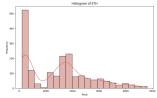
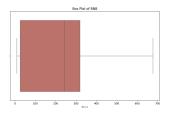


FIGURE 4. Ethereum price's histogram



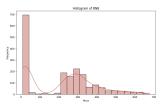


FIGURE 5. BNB price's boxplot

FIGURE 6. BNB price's histogram

IV. METHODOLOGY

A. LINEAR REGRESSION

Linear regression is an algorithm that provides a linear relationship between an independent variable and a dependent variable. A linear regression model with more than one independent variable is called a multiple linear regression model. 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.
- $X_1, ..., X_k$ 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.

B. AUTOREGRESSIVE INTEGRATED MOVING AVERAGE (ARIMA)

ARIMA is an acronym for AutoRegressive Integrated Moving Average. It is a statistical analysis model that uses time series data to either better understand the data set or to predict future trends. The full model can be written as:

$$y_t'=c+\phi_1y_{t-1}'+\ldots+\phi_py_{t-p}'+\theta_1\varepsilon_{t-1}+\ldots+\theta_q\varepsilon_{t-q}+\varepsilon_t$$
 Where:

- y'_t is the value of the time series at time t.
- c is the constant term.
- $\phi_1, ..., \phi_p$ are the autoregressive coefficients.
- $\theta_1, ..., \theta_q$ are the moving average coefficients.
- ε_t is the white noise at time t.

A non - seasonal ARIMA model is classified as an ARIMA (p, d, q) model, where:

- p is the number of lag observations included in the
- d is the number of times that the raw observations are differenced.
- q is the size of the moving average window, also called the order of moving average.

C. AUTOREGRESSIVE INTEGRATED MOVING AVERAGE WITH EXOGENOUS FACTORS (SARIMAX)

SARIMAX is an acronym for Seasonal Autoregressive Integrated Moving Average with eXogenous factors. The SARI-MAX model is an improved version of the SARIMA model, with exogenous factors (X) as external feature parameters for enhancing the model's performance, reducing the prediction errors, overcoming the autocorrelation issues, and improving

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the prediction results. The SARIMAX model consists of both seasonal effects and exogenous factors that can be used as SARIMAX (p, d, q) * (P, D, Q), while the exogenous factors are optional parameters. The exogenous factors are used to support the prediction model and to provide it with more details. The SARIMAX model can be presented as:

$$d_t = c + \sum_{n=1}^p \alpha_n d_{t-n} + \sum_{n=1}^q \theta_n \epsilon_{t-n} + \sum_{n=1}^r \beta_n x_{n_t} + \sum_{n=1}^p \phi_n d_{t-sn} + \sum_{n=1}^q \eta_n \epsilon_{t-sn} + \epsilon_t$$

Where:

- *P* is seasonal autoregressive order.
- D is seasonal differencing order.
- Q is seasonal moving average order.
- s is the length of the seasonal cycle.

D. RANDOM FOREST

Random forest is an ensemble learning method for classification, regression, and other tasks, that operates by constructing a multitude of decision trees at training time and outputting the class that is the mode of the classes (classification) or mean prediction (regression) of the individual trees.

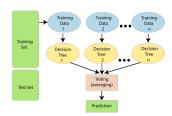


FIGURE 7. The structure of Random Forest model

Each tree in a random forest randomly samples subsets of the training data in a process known as bootstrap aggregating (bagging). The model is fit to these smaller data sets and the predictions are aggregated. Several instances of the same data can be used repeatedly through replacement sampling, and the result is that trees that are not only trained on different sets of data, but also different features used to make decisions.

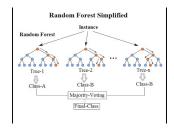


FIGURE 8. Overview of Random Forest model

The following steps explain the working Random Forest algorithm:

1) Select random samples from a given data or training set.

- This algorithm will construct a decision tree for every training data.
- 3) Voting will take place by averaging the decision tree.
- 4) For classification, use majority voting; for regression, use averaging to derive the final output.

E. RECURRENT NEURAL NETWORK (RNN)

Recurrent Neural Networks (RNNs) are a class of artificial neural networks specifically designed for processing sequences of data. Unlike traditional feedforward neural networks, RNNs have cyclical connections, allowing them to maintain a memory of previous inputs.

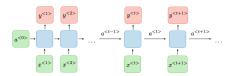


FIGURE 9. Overview of Recurrent Neural Network

For each timestep t, the activation $a^{< t>}$ and the output $y^{< t>}$ are expressed as follows:

$$a^{< t>} = g_1(W_{aa}a^{< t-1>} + W_{ax}x^{< t>} + b_a)$$

and

$$y^{< t>} = g_2(W_{ya}a^{< t>} + b_y)$$

Where:

- $W_{ax}, W_{as}, W_{ya}, b_a, b_y$ are coefficients that are shared temporally.
- g_1, g_2 are activation functions.

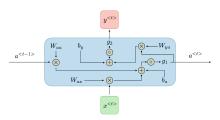


FIGURE 10. Architecture of a traditional RNN

In the case of a recurrent neural network, the loss function \mathcal{L} of all time steps is defined based on the loss at every time step:

$$\mathcal{L}(\hat{y},y) = \sum_{t=1}^{T_y} \mathcal{L}(\hat{y}^{< t>}, y^{< t>})$$
 Backpropagation is done at each point in time. At timestep T ,

Backpropagation is done at each point in time. At timestep T, the derivative of the loss \mathcal{L} with respect to weight matrix W is expressed as follows:

$$\left. \frac{\partial \mathcal{L}^{(T)}}{\partial W} = \sum_{t=1}^{T} \left. \frac{\partial \mathcal{L}^{(T)}}{\partial W} \right|_{(t)}$$

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