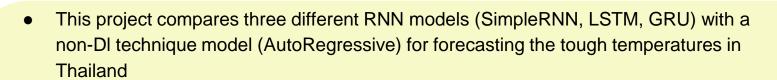


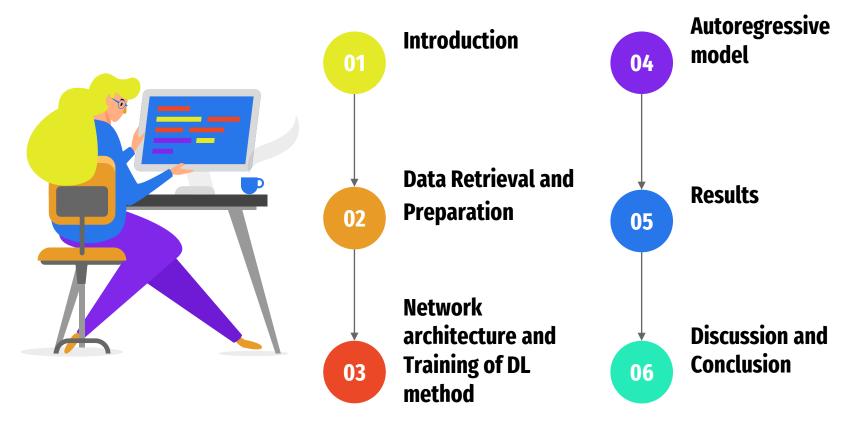
Highlights





- This study used a kaggle dataset to analyze how average temperatures in Thailand changed rapidly during Jan 1863 - Aug 2013.
- In terms of MAE comparison, you will see what the performance would be for various results of RNN with a non-DI technique model (Auto-Regressive).
- In this investigation, the interesting finding was that GRU model was the most potent model.
- LSTM and GRU train much faster than SimpleRNN

Agenda



Introduction

Some argue that global warming is the greatest peril our generation, while others argue that it is a fiction based on dubious science. In this study, we provide some data so that you can make your own assumptions.

There's a lot of data cleaning and preparation that goes into putting together a long-term study of weather patterns, even more than with other data sets that Kaggle has covered. Technicians used mercury thermometers to collect early data, and any fluctuation in visit duration had an impact on measurements.



Introduction



Many weather stations were relocated in the 1940s due to airport building. Electronic thermometers, which are thought to have a cooling bias, were particularly popular.

Given this intricacy, a multitude of different organizations gather information on climate trends. NOAA's MLOST, NASA's GISTEMP, and the UK's HadCrut are the three most widely used atmospheric and oceanic temperature data sets.

We repackaged the data from a more recent Berkeley Earth compilation, which is linked with Lawrence Berkeley National Laboratory. The Berkeley Earth Surface Temperature Study brings together 1.6 billion temperature reports from 16 repositories.

Introduction



It comes in a neat bundle and can be sliced into interesting subgroups (for example by country). They make the underlying data as well as the code for the alterations they performed public. They also employ approaches that allow for the inclusion of weather records from shorter time series, resulting in fewer observations being discarded.

The average land temperature in this dataset from Kaggle begins in 1750, while the maximum and minimum land temperatures, as well as worldwide ocean and land temperatures, commence in 1850.

Since we live in Thailand, we chose Thailand to focus on the small scale for this scope of study.

Purpose



The aims of this study are:

- To test the accuracy of average Thailand's temperature forecasts using four alternative models. *
- To compare the efficacy of four different model forecasts for average Thailand's temperatures.

* The four models in this study are:

1) RNN 2) LSTM 3) GRU 4) AutoRegressive

Data Retrieval

Climate Change: Earth Surface Temperature Data

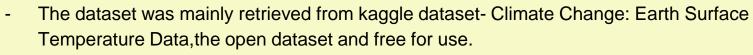
Exploring global temperatures since 1750



Data Code (558) Discussion (9) Metadata

About Dataset

Usability © 7.65

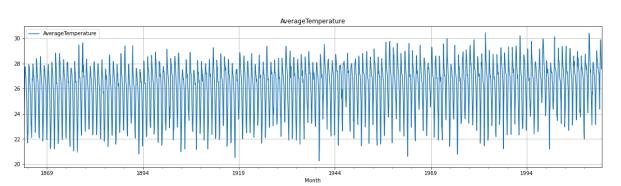


URL: https://www.kaggle.com/datasets/berkeleyearth/climate-change-earth-surface-temperature-data

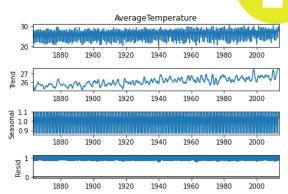
- "GlobalLandTemperaturesByCountry" was a csv file to employ in this study.
- There were 1808 rows x 4 columns during Jan 1863 Aug 2013.



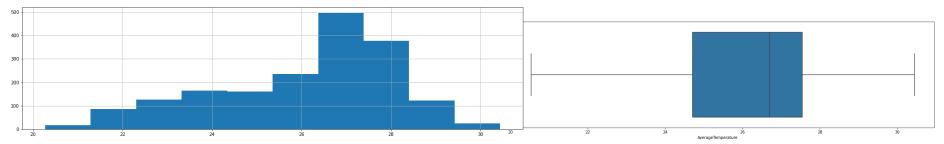
Data Exploration



Plot time series data



Multiplicative seasonal decomposition



Histogram plot

Box plot and interquartile range

Data Preparation

	dt	AverageTemperature	${\bf Average Temperature Uncertainty}$	Country
0	1743-11-01	4.384	2.294	Åland
1	1743-12-01	NaN	NaN	Åland
2	1744-01-01	NaN	NaN	Åland
3	1744-02-01	NaN	NaN	Åland
4	1744-03-01	NaN	NaN	Åland
577457	2013-05-01	19.059	1.022	Zimbabwe
577458	2013-06-01	17.613	0.473	Zimbabwe
577459	2013-07-01	17.000	0.453	Zimbabwe
577460	2013-08-01	19.759	0.717	Zimbabwe
577461	2013-09-01	NaN	NaN	Zimbabwe

577462 rows x 4 columns

Unique country names in the file: 243

Original data

	dt	AverageTemperature	AverageTemperatureUncertainty	Country	
0	1816-03-01	25.959	1.751	Thailand	
1	1816-04-01	27.263	2.960	Thailand	
2	1816-05-01	27.932	1.923	Thailand	
3	1816-06-01	26.500	1.940	Thailand	
4	1816-07-01	25.092	1.605	Thailand	
2366	2013-05-01	29.548	0.286	Thailand	
2367	2013-06-01	28.325	0.207	Thailand	
2368	2013-07-01	27.564	0.318	Thailand	
2369	2013-08-01	27.548	0.289	Thailand	
2370	2013-09-01	NaN	NaN	Thailand	
2371 rows × 4 columns					
#Row contain NaN data in each colunm dt 0					
AverageTemperature 125 AverageTemperatureUncertainty 125 Country 0					

Thailand data

- The original file contains temperature data for 243 unique countries
- Since our project focuses only on Thailand, the other countries data were filtered out using pandas

dtype: int64

However Thailand data still contains 125 rows of NaN data

Data Preparation

0 1863-01-01

2 1863-03-01

1807 2013-08-01

1863-02-01



Thailand

2.022 Thailand

0.854 Thailand

0.289 Thailand

	dt	AverageTemperature	$\label{lem:lemperature} \textbf{AverageTemperatureUncertainty}$	Country
46	1820-01-01	NaN	NaN	Thailand
93	1823-12-01	NaN	NaN	Thailand
94	1824-01-01	NaN	NaN	Thailand
95	1824-02-01	NaN	NaN	Thailand
96	1824-03-01	NaN	NaN	Thailand
558	1862-09-01	NaN	NaN	Thailand
559	1862-10-01	NaN	NaN	Thailand
560	1862-11-01	NaN	NaN	Thailand
561	1862-12-01	NaN	NaN	Thailand
2370	2013-09-01	NaN	NaN	Thailand

_				
3	1863-04-01	27.646	1.523	Thailand
4	1863-05-01	27.756	1.296	Thailand
1803	2013-04-01	29.885	0.234	Thailand
1804	2013-05-01	29.548	0.286	Thailand
1805	2013-06-01	28.325	0.207	Thailand
1806	2013-07-01	27.564	0.318	Thailand

22.806

24,700

26.599

Thailand NaN data

Thailand data after cleaning

27.548

- After exploring the NaN data, we found that the data after index 561 (dt=1862-12-01) did not contain any NaN data except the last row data
- So the data used will be the data from index 562 to 2369 (the row before the last row)

Data Preparation Code



```
df = pd.read csv('archive/GlobalLandTemperaturesByCountry.csv')
display(df)
\#https://www.kaqqle.com/datasets/berkeleyearth/climate-change-earth/-surface-temperature-data?select=GlobalTemperatures.csv
#LandAverageTemperature: global average land temperature in celsius
print('Unique country names in the file:', df['Country'].nunique())
df = df[df['Country']=='Thailand']
df = df.reset index(drop=True)
display(df)
print('#Row contain NaN data in each colunm')
print(df.isnull().sum())
display(df[df['AverageTemperature'].isnull()])
display(df[df['AverageTemperatureUncertainty'].isnull()])
#^uncomment the code above to see NaN data start from index 46 (1820-01-01) to 561 (1862-12-01) and 2370 (2013-09-01)
df = df[562:-1] #so we use data from index 562 to 2369
df = df.reset index(drop=True)
print(df.isnull().sum())
df
```

Data Preparation



```
column_data = df['AverageTemperature']

def convertToMatrix(data, feature_timestep):
    x, y = [], []
    for i in range(len(df)-feature_timestep):
        d = i+feature_timestep
        x.append(data[i:d])
        y.append(data[d])
    return np.array(x), np.array(y)
```

CODE: Transform raw data to training data set form

- Raw data will be transformed to deep learning desired form by calling **convertToMatrix** function
- Transformed shape depends on the number of feature timestep parameter called during each experiment

Data Preparation

```
feature timestep = 4
x, y = convertToMatrix(column data, feature timestep)
print(f'{feature_timestep} Features')
print(f'Data Length for {feature timestep} Features: {len(x)}')
print(f'Length x=y: \{len(x)==len(y)\}')
print(f'{feature timestep} Features x shape: {x.shape}')
print(f'{feature timestep} Features y shape: {y.shape}')
                                                                  percent trainset = 0.8
                                                                  percent testset = 0.1
4 Features
                                                                  rng = np.random.RandomState(random state seed)
Data Length for 4 Features: 1804
Length x=y: True
                                                                  rna.shuffle(x)
                                                                  rng = np.random.RandomState(random_state_seed)
4 Features x shape: (1804, 4)
                                                                  rng.shuffle(y)
4 Features y shape: (1804,)
                                                                  n train = int(len(x)*percent trainset)
                                                                  n \text{ test} = int(len(x)*percent testset)
                                                                  n \text{ val} = \text{len}(x)-n \text{ train-n test}
                                                                  x train, y train = x[:n train], y[:n train]
                                                                  x val, y val = x[n train:n train+n val], y[n train:n train+n val]
                                                                  x test, y test = x[n train+n val:], y[n train+n val:]
                                                                  print('Split Train/Val/Test')
                                                                  print(f'{feature timestep} Features Trainset Shape:',x train.shape, y train.shape)
                                                                  print(f'{feature timestep} Features Valset Shape:',x val.shape, y val.shape)
                                                                  print(f'{feature timestep} Features Testset Shape:',x test.shape, y test.shape)
                                                                  Split Train/Val/Test
                                                                  4 Features Trainset Shape: (1443, 4) (1443,)
```

4 Features Valset Shape: (181, 4) (181,) 4 Features Testset Shape: (180, 4) (180,)

Data Preparation: Normalized all data by x_train

```
#normalized all data by x train
minmax norm = MinMaxScaler().fit(x train.reshape(-1,1))
x train_norm = minmax_norm.transform(x_train.reshape(-1,1)).reshape(-1,feature_timestep)
x val norm = minmax norm,transform(x val,reshape(-1,1)),reshape(-1,feature timestep)
x test norm = minmax norm.transform(x test_reshape(-1,1)),reshape(-1,feature timestep)
y train norm = minmax norm.transform(y train.reshape(-1,1)).reshape(-1,1)
y_{val} norm = minmax_norm.transform(y_{val}.reshape(-1,1)).reshape(-1,1)
v test norm = minmax norm.transform(v test.reshape(-1.1)).reshape(-1.1)
print('x shape before newaxis')
                                                              print('x shape after newaxis')
print(x train norm.shape)
                                                              print(x train norm.shape) #Final input shape must be (n sample, n sequence, n feature per sequence) https://www.voutube.com/watch?v=EnuAP1ZOb4s
print(x val norm.shape)
                                                              print(x val norm.shape)
print(x_test_norm.shape)
                                                              print(x test norm.shape)
                                                              print('y shape')
 #add new axis
                                                              print(v train norm.shape)
x train norm = x train norm[..., np.newaxis]
 "need input as [[],
                                                              print(y val norm.shape)
                                                              print(y test norm.shape)
                                                               #transform to float32
            []] shape=(4,1)
                                                              x train norm = x train norm.astype(np.float32)
                                                              x val norm = x val norm.astype(np.float32)
x_{val_norm} = x_{val_norm}[..., np.newaxis]
x test norm = x_test_norm[..., np.newaxis]
                                                              x test norm = x test norm.astype(np.float32)
                                                              v train norm = v train norm.astvpe(np.float32)
                                                              v val norm = v val norm.astvpe(np.float32)
                                                              y test norm = y test norm.astype(np.float32)
                                                              x shape before newaxis
                                                              (1443, 4)
                                                               (181, 4)
                                                               (180, 4)
                                                              x shape after newaxis
                                                              (1443, 4, 1)
                                                               (181, 4, 1)
                                                              (180, 4, 1)
                                                              y shape
                                                               (1443, 1)
                                                               (181, 1)
                                                               (180, 1)
```



Network Architecture: Deep RNN

- Deep RNN model consists of
 - ☐ SimpleRNN layer (128 neurons, activation=tanh)
 - ☐ **Dropout** layer (**0.5** dropout rate)
 - SimpleRNN layer (128 neurons, activation=tanh)
 - ☐ **Dropout** layer (**0.5** dropout rate)
 - □ Dense layer (64 neurons)
 - ☐ **Dropout** layer (**0.5** dropout rate)
 - □ Dense layer (1 neuron)

```
#deep rnn model
model = Sequential()

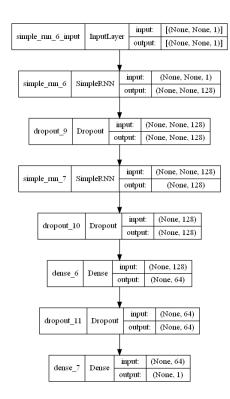
model.add(SimpleRNN(128, input_shape=(None,1), return_sequences=True))
model.add(Dropout(0.5))

model.add(SimpleRNN(128))
model.add(Dropout(0.5))

model.add(Dense(64))
model.add(Dropout(0.5))

model.add(Dense(1))
```

CODE: Deep RNN model



Deep RNN model structure

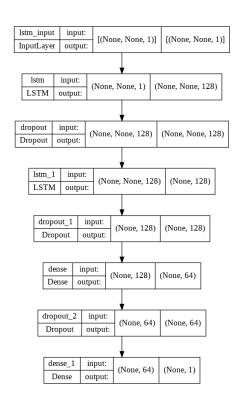


Network Architecture: LSTM

- Deep RNN model consists of
 - ☐ LSTM layer (128 neurons, activation=tanh)
 - ☐ **Dropout** layer (**0.5** dropout rate)
 - ☐ LSTM layer (128 neurons, activation=tanh)
 - ☐ **Dropout** layer (**0.5** dropout rate)
 - Dense layer (64 neurons)
 - □ **Dropout** layer (**0.5** dropout rate)
 - □ Dense layer (1 neuron)

```
1 #deep LSTM model
2 model = Sequential()
3
4 model.add(LSTM(128, input_shape=(None,1), return_sequences=True))
5 model.add(Dropout(0.5))
6
7 model.add(LSTM(128))
8 model.add(Dropout(0.5))
9
10 model.add(Dense(64))
11 model.add(Dropout(0.5))
12
13 model.add(Dense(1))
```

CODE: LSTM model



LSTM model structure



Network Architecture: GRU

- Deep RNN model consists of
 - ☐ GRU layer (128 neurons, activation=tanh)
 - ☐ **Dropout** layer (**0.5** dropout rate)
 - □ GRU layer (128 neurons, activation=tanh)
 - ☐ **Dropout** layer (**0.5** dropout rate)
 - □ Dense layer (64 neurons)
 - □ **Dropout** layer (**0.5** dropout rate)
 - □ Dense layer (1 neuron)

```
# GRU model
model = Sequential()

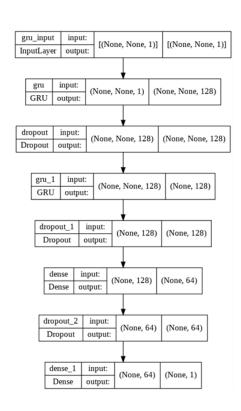
model.add( GRU(128, input_shape=(None,1), return_sequences=True))
model.add(Dropout(0.5))

model.add(GRU(128))
model.add(Dropout(0.5))

model.add(Dense(64))
model.add(Dropout(0.5))

model.add(Dense(1))
```

CODE: GRU model



GRU model structure



 All deep learning models (Deep RNN, LSTM and GRU) the training/validation//testing dataset are as follows:

```
training set:
    x_train_norm, y_train_norm

validation set:
    x_val_norm, y_val_norm

testing set:
    x_test_norm
```

 All deep learning models (Deep RNN, LSTM and GRU) the training parameters was set as follows:

training parameters:

optimizer = adam(learning rate=0.001, decay=1e-6)

loss = mean absolute error

(MAE was selected instead of RMSE because the error calculated from predicted output is a normalized output which is a decimal number, using RMSE will lower down an actual error, so MAE was selected to truly represent an actual error)

performance metric = mean absolute error

The best trained model was saved using **ModelCheckpoint** monitored on **minimizing** validation loss

epochs = 100

batch size = 10

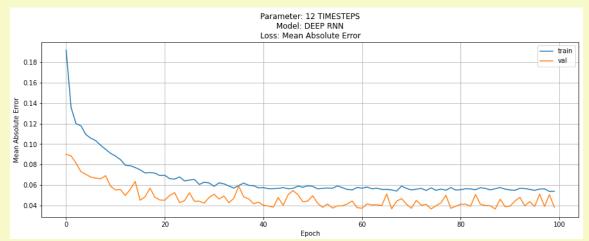


CODE: Training

Model: "sequential_8"				
Layer (type)	Output Shape	Param #		
simple_rnn_16 (SimpleRNN)		16640		
dropout_24 (Dropout)	(None, None, 128)	0		
simple_rnn_17 (SimpleRNN)	(None, 128)	32896		
dropout_25 (Dropout)	(None, 128)	0		
dense_16 (Dense)	(None, 64)	8256		
dropout_26 (Dropout)	(None, 64)	0		
dense_17 (Dense)	(None, 1)	65		
Total params: 57,857 Trainable params: 57,857 Non-trainable params: 0				
<pre>C:\Users\tOm\AppData\Roamin precated, use `learning_rat super(Adam, self)init_</pre>	e` instead.	packages\keras\optimi	mizer_v2\adam.py:105: UserWarning: The `lr` argument i	s de
144/144 [_mean_absolute_error: 0.193 Epoch 2/100		ms/step - loss: 0.782	825 - mean_absolute_error: 0.7825 - val_loss: 0.1937 -	val
144/144 [====== _mean_absolute_error: 0.079		ms/step - loss: 0.338	380 - mean_absolute_error: 0.3380 - val_loss: 0.0793 -	val
_mean_absolute_error: 0.086		ms/step - loss: 0.214	144 - mean_absolute_error: 0.2144 - val_loss: 0.0868 -	val
Epoch 4/100 144/144 [_mean_absolute_error: 0.058		ms/step - loss: 0.167	676 - mean_absolute_error: 0.1676 - val_loss: 0.0586 -	val
Epoch 5/100 144/144 [mean absolute error: 0.057		ms/step - loss: 0.141	416 - mean_absolute_error: 0.1416 - val_loss: 0.0572 -	val
Epoch 6/100	=======] - 16s 112r	ms/step - loss: 0.120	209 - mean_absolute_error: 0.1209 - val_loss: 0.0617 -	val

Training output

For each model, the plot of MAE in training and validation set on each training epoch is
plotted see how the training goes and to detect if there was any abnormal error during the
training



Training set and validation set MAE on each training epoch (Example from Deep RNN with 12 timesteps feature)

```
#plot loss, val_loss VS epoch
loss_metric = f'Mean Absolute Error'
plt.figure(figsize=(15,5))
plt.plot(history.history['loss'])
plt.plot(history.history['val_loss'])
plt.title(f'Parameter: {feature_timestep} TIMESTEPS\n\
Model: {model_structure}\nLoss: {loss_metric}')
plt.ylabel(f'{loss_metric}')
plt.xlabel('Epoch')
plt.legend(['train','val'], loc='upper right')
plt.grid()
plt.show()
```

CODE: MAE plotting

- Then the best model from training session was loaded
- The best model was tested on validation set to see how the model perform (the loss/metric from this session still represents in normalized value)

CODE: Loading best model and validating on validation set (Example from Deep RNN with 12 timesteps feature)

- The actual prediction for test set was performed by the best model and then inverse transform(denormalized) to present an actual predicted temperature
- The results will be further interpreted and concluded in result section

```
y_test_predict = model.predict(x_test_norm)
y_test_predict_inv = minmax_norm.inverse_transform(y_test_predict)
```

CODE: Test set temperature prediction and denormalization

 The deep learning model training time comparison (from command %%time) with machine specifications as followed are shown in the table below

Machine Specifications:

CPU: Intel(R) Core(TM) i5-8300H CPU @ 2.30GHz 2.30 GHz

o RAM: 24.0 GB (23.9 GB usable)

o GPU: NVIDIA GeForce GTX 1050Ti

Model	Timesteps Feature			
	4	6	12	
Deep RNN	11m 28s	16m 6s	27m 54s	
LSTM	3m 59s	3m 55s	4m 19s	
GRU	3m 54s	3m 38s	3m 35s	

Auto Regressive model

In order to deal with the forecasting task of time-series dataset, auto regressive model is usually chosen to be the one to predict the outcome of the future values based on previous data values.

$$y_t = c + \phi_1 y_{t-1} + \phi_2 y_{t-2} + ... + \phi_p y_{t-p} + \epsilon_t$$

An autoregressive model uses a linear combination of past values of the target to make forecasts and is made against the target itself. Where lags express each data point across time is called a lag, Bias and weight are associated with each lag which tells the importance of that time step in predicting the final value and Autocorrelation is the relationship between forecasted variable and input variables.

Finding autocorrelation

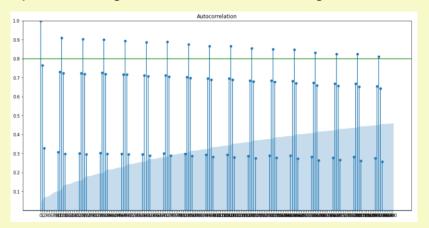


In order to find out which lags have a strong correlation to the forecasted values, The code creates a plot that shows us how much each previous lag influences the future lag.

```
# find out which is the best lag to use
import pandas as pd
import matplotlib.pyplot as plt
from statsmodels.graphics.tsaplots import plot_acf
import numpy as np

fig, ax = plt.subplots(figsize=(16,8))
plot_acf(df1['AverageTemperature'], lags=200, ax=ax)
plt.ylim([0,1])
plt.yticks(np.arange(0.1, 1.1, 0.1))
plt.xticks(np.arange(0, 201, 1))
plt.axhline(y=0.8, color="green")
plt.show()
```

CODE: Finding which lag is the best lag to use



Autocorrelation and lag plot

Then we provide a threshold of 0.8 to see the lag that has a very strong correlation to the future value. The lag that we are looking for is 192 which goes above the threshold.

Training Auto Regressive model

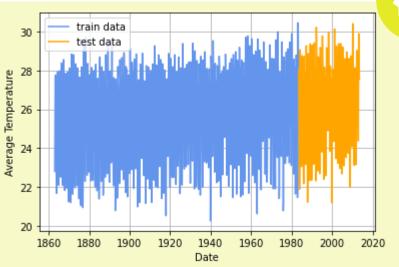
```
train_df = df1.iloc[:1446]
test_df = df1.iloc[:1446]
print('shapes:', df1.shape, train_df.shape, test_df.shape)

shapes: (1808, 1) (1446, 1) (362, 1)

train_set_size = len(train_df)
train_set_dates = df1.head(train_set_size).index # for plotting
test_set_dates = df1.tail(362).index

plt.grid()
plt.plot(train_set_dates, train_df.AverageTemperature, color='cornflowerblue', label='train data')
plt.plot(test_set_dates, test_df.AverageTemperature, color='orange', label='test data')
plt.legend(loc='best')
plt.ylabel('bate')
plt.ylabel('bate')
plt.show()
```

CODE: Splitting training set and testing set



Date and time that used for training and testing

```
from statsmodels.tsa.ar_model import AutoReg

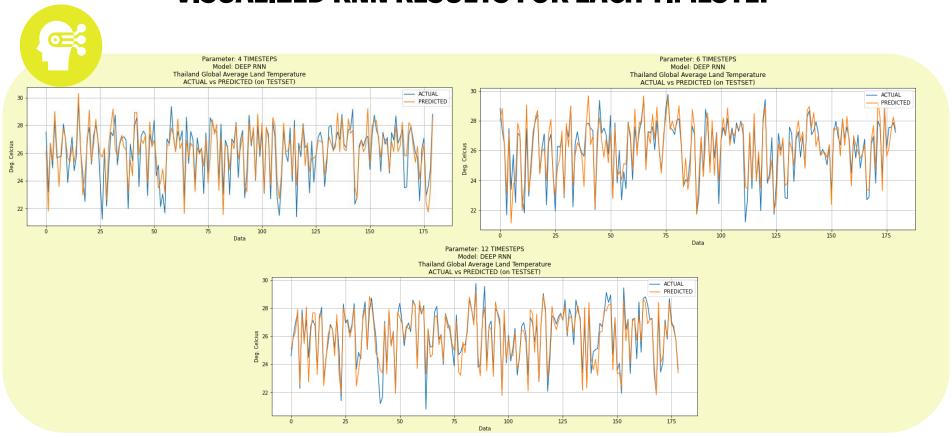
model = AutoReg(train_df, lags=192)
trained_model = model.fit()
```

CODE: Creating and fitting the model

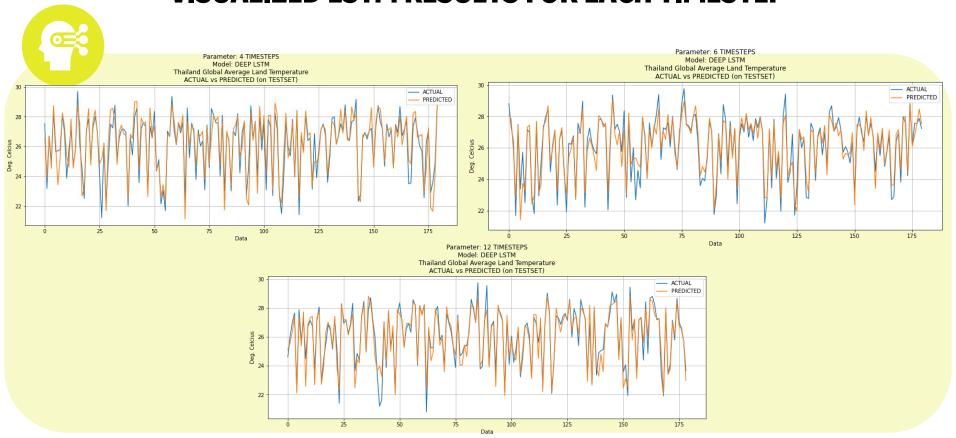


Results

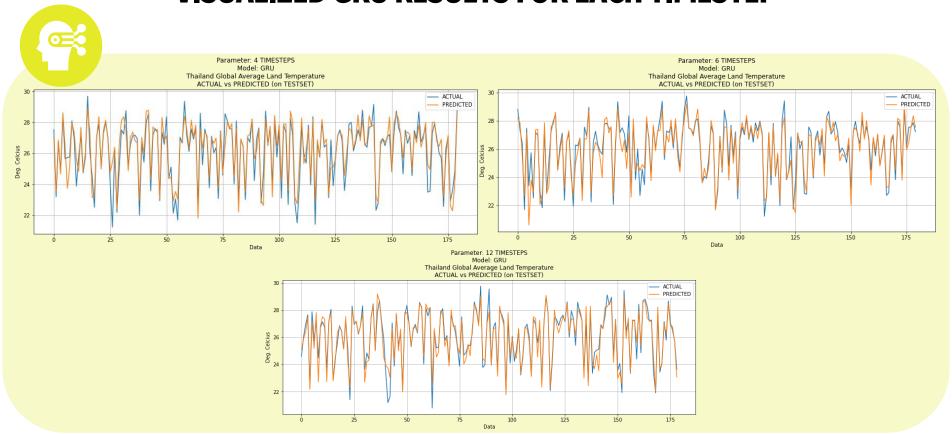
VISUALIZED RNN RESULTS FOR EACH TIMESTEP



VISUALIZED LSTM RESULTS FOR EACH TIMESTEP

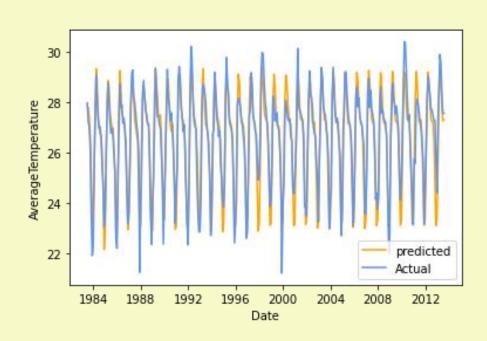


VISUALIZED GRU RESULTS FOR EACH TIMESTEP



VISUALIZED AUTOREGRESSIVE RESULTS





MAE Evaluation on Test Set



Models	4 Times	6 Times	12 Times
RNN	0.785	0.650	0.516
LTSM	0.576	0.493	0.433
GRU	0.551	0.482	0.411
AutoRegressive	ve 0.531		

Conclusion



According to the experiment, the all deep learning's best models (Multi-layer RNN, LSTM and GRU) have better performance compared to the autoregressive method. Comparing among the deep learning model, the GRU model has minimum MAE at 0.411. The second best is LSTM at 0.433 and the last on is RNN at 0.516. We'll find that the GRU and LSTM model outperform the RNN model.

In term of input time step which is one of the important hyper-parameters, we found that the 12 time-steps input made the model the best outcome comparing to 4 and 6 time-steps input. There is a point to mentioned that the data were collected on the monthly basis and the temperature data is affected from the season of the year which may be the reason that 12 time-steps input show the best result in these experiments.

Discussion

- The results are quite good with MAE around 0.4 0.8 which is acceptable for temperature prediction. However, in some cases which require more accurate prediction, the configuration of GRU and LSTM model should be considered to achieve more accurate model.
- For other cases which the data may not be collected on the monthly basis, the best input time step may not be 12 as per these experiments. The input time step should be taken into account and properly fine-tuned based on each data.

References



Python Version: 3.8.5

Python Library

- Matplotlib 3.5.1
- Numpy 1.22.0
- OpenCV 4.5.5
- Pandas 1.3.0
- Tensorflow 2.7.0

Source Code

 Data preparation and model training code referenced from RNN_ex1.py (BADS7604 by Asst. Prof. Thitirat Sribiborbornratanakul)

Datasets

• From Climate Change: Earth Surface Temperature Data
(https://www.kaggle.com/datasets/berkeleyearth/climate-change-earth-surface-temperature-data?select=GlobalTemperatures.csv) with license: Attribution-NonCommercial-ShareAlike 4.0 International (CC BY-NC-SA 4.0)

Citing



For those who want to use any presentation images please reference the image in **bibtex** format

Team Members

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(25%) Highlights, Data preparation, Network Architecture, Experimental variation on LSTM, Conclusion and Discussion

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Introduction and Highlights composition, Data preparation, Network Architecture, Experimental variation on GRU mode

End Credit



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Thank you for your attention