Time Series Foundamentals

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1 Topic: Mean Squared Prediction

Suppose $\{X_t\}$ is a stationary process with mean μ , and auto-covariance function $\gamma_X(h)$. Our goal is to predict: X_{n+h} using observations X_1, X_2, \ldots, X_n .

We define the Best Linear Predictor to be the one minimizes the Mean Squared Prediction Error (MSPE):

$$\widehat{X}_{n+h} = E\left[X_{n+h} \mid X_1, \dots, X_n\right] = P_n X_{n+h} := a_0^* + a_1^* X_n + \dots + a_n^* X_1 = a_0 + \mathbf{a}_n^\top \mathbf{x}_n = \arg\min_{a_0, \mathbf{a}_n} \mathbb{E}\left[\left(X_{n+h} - a_0 - \mathbf{a}_n^\top \mathbf{x}_n\right)^2\right] = \arg\min_{a_0, \mathbf{a}_n} L(a_0, \mathbf{a}_n)$$

where $\mathbf{a}_n=(a_1,\ldots,a_n)\in\mathbb{R}^n$, $\mathbf{x}_n=(x_n,\ldots,x_1)\in\mathbb{R}^n$, P in P_n stands for "Predictor" and n suggests that we are using n most recent observations.

To simplify our computation, we define
$$\mathbb{1}=(1,\dots,1)\in\mathbb{R}^n$$
, $\boldsymbol{\gamma}_n(h)=\begin{bmatrix} \gamma_X(h)\\ \gamma_X(h+1)\\ \vdots\\ \gamma_X(h+n-1)\end{bmatrix}\in\mathbb{R}^n$, $\boldsymbol{\Gamma}_n=\begin{bmatrix} \gamma_X(0)&\gamma_X(1)&\cdots&\gamma_X(n-1)\\ \gamma_X(1)&\gamma_X(0)&\cdots&\gamma_X(n-2)\\ \vdots&\vdots&\ddots&\vdots\\ \gamma_X(n-1)&\gamma_X(n-2)&\cdots&\gamma_X(0)\end{bmatrix}\in\mathbb{R}^{n\times n}$.

Taking the partial derivatives of L, we obtain

$$0 = \frac{\partial L}{\partial a_0} = \mathbb{E}\left[\frac{\partial}{\partial a_0}(X_{n+h} - a_0 - \mathbf{a}_n^{\top}\mathbf{x}_n)^2\right] = \mathbb{E}\left[-2(X_{n+h} - a_0 - \mathbf{a}_n^{\top}\mathbf{x}_n)\right] = 2(a_0 - \mu - \mathbf{a}_n^{\top}\mathbb{1}) \Longrightarrow a_0^* = \mu(1 - \mathbf{a}_n^{\top}\mathbb{1})$$

and

$$\begin{aligned} \mathbf{0} &= \frac{\partial L}{\partial \mathbf{a}_n} = \frac{\partial}{\partial \mathbf{a}_n} \mathbb{E} \left[(X_{n+h} - a_0 - \mathbf{a}_n^\top \mathbf{x}_n)^2 \right] \\ &= \mathbb{E} \left[\frac{\partial}{\partial \mathbf{a}_n} \left(\mathbf{a}_n^\top \mathbf{x}_n \mathbf{x}_n^\top \mathbf{a}_n - 2X_{n+h} \mathbf{x}_n^\top \mathbf{a}_n + 2a_0 \mathbf{x}_n^\top \mathbf{a}_n + (\text{ terms without } \mathbf{a}_n) \right) \right] \\ &= 2\mathbb{E} [\mathbf{x}_n \mathbf{x}_n^\top \mathbf{a}_n - X_{n+h} \mathbf{x}_n + a_0 \mathbf{x}_n] \\ &= 2\mathbb{E} [\mathbf{x}_n \mathbf{x}_n^\top] \mathbf{a}_n - 2\mathbb{E} [X_{n+h} \mathbf{x}_n] + 2\mathbb{E} [a_0 \mathbf{x}_n] \\ &= 2(\Gamma_n + \mu^2 \mathbb{1} \mathbb{1}^\top) \mathbf{a}_n - 2(\gamma_n(h) + \mu^2 \mathbb{1}) + 2a_0 \mu \mathbb{1} \\ a_0 &= \mu (1 - \mathbf{a}_n^\top \mathbb{1}) \Longrightarrow = 2(\Gamma_n + \mu^2 \mathbb{1} \mathbb{1}^\top) \mathbf{a}_n - 2(\gamma_n(h) + \mu^2 \mathbb{1}) + 2\mu (1 - \mathbf{a}_n^\top \mathbb{1}) \mu \mathbb{1} \\ &= 2\Gamma_n \mathbf{a}_n - 2\gamma_n(h) \\ \Longrightarrow \mathbf{a}_n^* &= \Gamma_n^{-1} \gamma_n(h) \end{aligned}$$

Therefore, the Best Linear Predictor is given by

$$P_n X_{n+h} := a_0^* + \mathbf{a}_n^{*^\top} \mathbf{x}_n \text{ where } \begin{cases} a_0^* &= \mu (1 - \mathbf{a}_n^{*^\top} \mathbb{1}) \\ \mathbf{a}_n^* &= \Gamma_n^{-1} \gamma_n(h) \end{cases}$$
(1)

Problem 1.1. Show that Mean Squared Predictive Error is obtained by

$$\mathbb{E}\left[\left(X_{n+h} - P_n X_{n+h}\right)^2\right] = \gamma_X(0) - \mathbf{a}_n^{*^{\top}} \boldsymbol{\gamma}_n(h)$$

Solution: From above, we get

$$\begin{split} \mathbb{E}\left[\left(X_{n+h}-P_{n}X_{n+h}\right)^{2}\right] &= \mathbb{E}\left[\left(X_{n+h}-a_{0}^{*}-\mathbf{a}_{n}^{*^{\top}}\mathbf{x}_{n}\right)^{2}\right] \\ \mathbf{a}_{n}^{*^{\top}}\mathbf{x}_{n} &= \mathbf{x}_{n}^{\top}\mathbf{a}_{n}^{*} \Longrightarrow = \mathbb{E}\left[X_{n+h}^{2}+a_{0}^{*^{2}}+\mathbf{a}_{n}^{*^{\top}}\mathbf{x}_{n}\mathbf{x}_{n}^{\top}\mathbf{a}_{n}^{*}-2a_{0}^{*}X_{n+h}-2X_{n+h}\mathbf{a}_{n}^{*^{\top}}\mathbf{x}_{n}+2a_{0}^{*}\mathbf{a}_{n}^{*^{\top}}\mathbf{x}_{n}\right] \\ &= \mathbb{E}\left[X_{n+h}^{2}\right]+a_{0}^{*^{2}}+\mathbf{a}_{n}^{*^{\top}}\mathbb{E}[\mathbf{x}_{n}\mathbf{x}_{n}^{\top}]\mathbf{a}_{n}^{*}-2a_{0}^{*}\mathbb{E}[X_{n+h}]-2\mathbf{a}_{n}^{*^{\top}}\mathbb{E}[X_{n+h}\mathbf{x}_{n}]+2a_{0}^{*}\mathbf{a}_{n}^{*^{\top}}\mathbb{E}[\mathbf{x}_{n}] \\ &= \gamma_{X}(0)+\mu^{2}+\mu^{2}(1-\mathbf{a}_{n}^{*^{\top}}\mathbf{1})^{2}+\mathbf{a}_{n}^{*^{\top}}(\Gamma_{n}+\mu^{2}\mathbf{1}\mathbf{1}^{\top})\mathbf{a}_{n}^{*}-2\mu^{2}(1-\mathbf{a}_{n}^{*^{\top}}\mathbf{1})-2\mathbf{a}_{n}^{*^{\top}}(\gamma_{n}(h)+\mu^{2}\mathbf{1})+2\mu^{2}(1-\mathbf{a}_{n}^{*^{\top}}\mathbf{1})\mathbf{a}_{n}^{*^{\top}}\mathbf{1} \\ &= \gamma_{X}(0)+\mathbf{a}_{n}^{*^{\top}}\Gamma_{n}\mathbf{a}_{n}^{*}-2\mathbf{a}_{n}^{*^{\top}}\gamma_{n}(h) \\ \mathbf{a}_{n}^{*}=\Gamma_{n}^{-1}\gamma_{n}(h) \Longrightarrow = \gamma_{X}(0)+\mathbf{a}_{n}^{*^{\top}}\Gamma_{n}\Gamma_{n}^{-1}\gamma_{n}(h)-2\mathbf{a}_{n}^{*^{\top}}\gamma_{n}(h) \\ &= \gamma_{X}(0)-\mathbf{a}_{n}^{*^{\top}}\gamma_{n}(h) \end{split}$$

Hence, we can conclude that the Mean Squared Predictive Error is obtained by

$$\mathbb{E}\left[\left(X_{n+h} - P_n X_{n+h}\right)^2\right] = \gamma_X(0) - \mathbf{a}_n^{*^{\top}} \boldsymbol{\gamma}_n(h)$$

Remark 1.2.

1. For a linear Gaussian process $\{X_t\}$, take $X_{n+h} \mid X_1, \dots, X_n = P_n X_{n+h} + \epsilon = a_0^* + \mathbf{a}_n^{*^\top} \mathbf{x}_n + \epsilon$, then the residual term ϵ must be Gaussian. Then, we have

$$\mathbb{E}[X_{n+h} \mid X_1, \dots, X_n - P_n X_{n+h}] = \mu - a_0^* + \mu \mathbf{a}_n^{*^{\top}} \mathbb{1} = 0 \xrightarrow{a_0^* = \mu(1 - \mathbf{a}_n^{*^{\top}} \mathbb{1})} \epsilon \sim \mathcal{N}(0, \text{MSPE}(P_n X_{n+h})) = \mathcal{N}(0, \gamma_X(0) - \mathbf{a}_n^{*^{\top}} \boldsymbol{\gamma}_n(h))$$

by our computation; thus, the predictive distribution is also Gaussian:

$$X_{n+h} \mid X_1, \dots, X_n \sim \mathcal{N}\left(P_n X_{n+h}, \text{MSPE}\left(P_n X_{n+h}\right)\right) = \mathcal{N}\left(a_0^* + \mathbf{a}_n^{*^{\top}} \mathbf{x}_n, \gamma_X(0) - \mathbf{a}_n^{*^{\top}} \boldsymbol{\gamma}_n(h)\right)$$

where $a_0^* = \mu(1 - \mathbf{a}_n^{*^{\top}} \mathbb{1})$ and $\mathbf{a}_n^* = \Gamma_n^{-1} \gamma_n(h)$.

2. For large n, Γ_n^{-1} is obtained through recursive algorithms. As an example look at Durbin-Levinson Algorithm

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2 Topic: MLE for AR and MA processes

Direct MLE

Suppose that $\{X_t\}$ is a Gaussian time series (process) with mean μ and autocovariance function $\gamma_X(i,j)$. Let $\mathbf{x}_n=(X_1,\ldots,X_n)\in\mathbb{R}^n$. Let Γ_n denote the covariance matrix $\Gamma_n=Cov\left(\mathbf{x}_n,\mathbf{x}_n\right)$, and assume that Γ_n is nonsingular. Since any subset of a Gaussian process are jointly Gaussian, the likelihood function for given observation \mathbf{x}_n (a single observation) is

$$L(\Psi) = f(\mathbf{x}_n \mid \Psi) = \frac{1}{\sqrt{(2\pi)^n |\Gamma_n|}} \exp\left(-\frac{1}{2}(\mathbf{x}_n - \mu)^\top \Gamma_n^{-1}(\mathbf{x}_n - \mu)\right).$$

where Ψ is the parameter set, e.g.for the ARMA model, $\Psi_{ARMA(p,q)} = (\phi_1, \dots, \phi_p, \theta_1, \dots, \theta_q, \sigma^2)$.

To simplify our computation, we could consider the log-likelihood function:

$$\log L(\Psi) = -\frac{n}{2}\log(2\pi) + \frac{1}{2}\log|\Gamma_n|^{-1} - \frac{1}{2}(\mathbf{x}_n - \boldsymbol{\mu})^{\top}\Gamma_n^{-1}(\mathbf{x}_n - \boldsymbol{\mu})$$

Then, the values of parameters Ψ which maximizes the Log-likelihood function is the Maximum Likelihood Estimate of Ψ . That is,

$$\hat{\Psi}_{MLE} = \arg\max_{\Psi} \log L(\Psi)$$

Conditional MLE

For the ease of implementation, we can also use the conditional MLE. Using the fact that

$$f(\mathbf{x}_n \mid \Psi) = f(x_1, \dots, x_n \mid \Psi) = f(x_n \mid x_1, \dots, x_{n-1}, \Psi) f(x_1, \dots, x_{n-1} \mid \Psi) \stackrel{\dots}{=} f(x_n \mid x_1, \dots, x_{n-1}, \Psi) f(x_{n-1} \mid x_1, \dots, x_{n-2}, \Psi) \cdots f(x_1 \mid \Psi)$$

Then, the MLE is given by

$$\log L(\Psi) = \log f(x_1 \mid \Psi) + \sum_{i=2}^{n} \log f(x_i \mid x_1, \dots, x_{i-1}, \Psi)$$
(2)

Similarly, the values of parameters Ψ which maximizes the Log-likelihood function is the Maximum Likelihood Estimate of Ψ . That is,

$$\hat{\Psi}_{MLE} = \arg\max_{\Psi} \log L(\Psi)$$

Remark 2.1.

- 1. We mostly use Conditional MLE in applications.
- 2. Since X_1 is the the first data point, the way it's picked is on its own. If X_1 is pick based on some distribution, we can surely keep the term $\log f(x_1 \mid \Psi)$ in the above equation and work with the exact MLE. If how X_1 is picked is unknown, we can treat it as deterministic, and drop the term $\log f(x_1 \mid \Psi)$ in the above equation.

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Problem 2.2. Using a programming language of your choice, repeat the MLE construction and simulation for AR and MA models as discussed this in the class.

Solution:

(a) Conditional MLE for AR(1) Process

Consider observations $\{X_1, X_2, \dots, X_n\}$ of a stationary Gaussian AR(1) process. For parameter vector is $\Psi = (\phi_0, \phi, \sigma^2)$, we have

$$X_t = \phi_0 + \phi X_{t-1} + Z_t, Z_t \sim IIDN\left(0,\sigma^2\right), \quad \mathbb{E}[X_t] = \frac{\phi_0}{1-\phi}, Var\left(X_t\right) = \frac{\sigma^2}{1-\phi^2} \text{ and } \gamma_X(h) = \frac{\phi^h \sigma^2}{1-\phi^2}$$

The probability distribution for X_1 can be written as,

$$X_1 \sim \mathcal{N}\left(\frac{\phi_0}{1 - \phi}, \frac{\sigma^2}{1 - \phi^2}\right) \Longrightarrow f\left(x_1; \Psi\right) = \frac{1}{\sqrt{2\pi \frac{\sigma^2}{1 - \phi^2}}} \exp\left\{\frac{-\left(x_1 - \frac{\phi_0}{1 - \phi}\right)^2}{2\frac{\sigma^2}{1 - \phi^2}}\right\}$$

Since $X_2 = \phi_0 + \phi X_1 + Z_2$, $X_2 \mid X_1 = x_1 \sim \mathcal{N}\left(\phi_0 + \phi x_1, \sigma^2\right)$. The conditional probability distribution for X_2 knowing X_1 can be written as.

$$f(x_2 \mid x_1; \Psi) = \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left\{ \frac{-(x_2 - \phi_0 - \phi x_1)^2}{2\sigma^2} \right\}$$

Similarly, we have $X_n \mid X_{n-1} = x_{n-1} \sim \mathcal{N}\left(\phi_0 + \phi x_{n-1}, \sigma^2\right)$ and

$$f(x_n \mid x_{n-1}; \Psi) = \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left\{ \frac{-(x_n - \phi_0 - \phi x_{n-1})^2}{2\sigma^2} \right\}$$

Plug these in Equation (2), we obtain the Log-Likelihood function:

$$\log L(\Psi) = \log f(x_1 \mid \Psi) + \sum_{i=2}^{n} \log f(x_i \mid x_1, \dots, x_{i-1}, \Psi)$$

$$= -\frac{1}{2} \log(2\pi) - \frac{1}{2} \log\left(\frac{\sigma^2}{1 - \phi^2}\right) - \frac{\left(x_1 - \frac{\phi_0}{1 - \phi}\right)^2}{2\frac{\sigma^2}{1 - \phi^2}} + \sum_{i=2}^{n} -\frac{1}{2} \log(2\pi) - \frac{1}{2} \log\left(\sigma^2\right) - \frac{\left(x_i - \phi_0 - \phi x_{i-1}\right)^2}{2\sigma^2}$$

$$= -\frac{1}{2} \log(2\pi) - \frac{1}{2} \log\left(\frac{\sigma^2}{1 - \phi^2}\right) - \frac{\left(x_1 - \frac{\phi_0}{1 - \phi}\right)^2}{2\frac{\sigma^2}{1 - \phi^2}} - \frac{n - 1}{2} \log(2\pi) - \frac{n - 1}{2} \log\left(\sigma^2\right) - \sum_{i=2}^{n} \frac{\left(x_i - \phi_0 - \phi x_{i-1}\right)^2}{2\sigma^2}$$

Maximize this function to estimate the parameters. (Much easier to implement)

Example: Suppose we have AR(1) process $X_t = 0.9X_{t-1} + Z_t, Z_t \sim IIDN (0, 0.7^2)$. We simulate this process and pretend that we do not know the true values for ϕ and σ . Then, we use the conditional MLE to estimate the parameter vector $\Psi = (\phi, \sigma^2)$ By the above deduction, we get the log likelihood function is

$$\log L(\Psi) = -\frac{1}{2}\log(2\pi) - \frac{1}{2}\log\left(\frac{\sigma^2}{1-\phi^2}\right) - \frac{x_1^2}{2\frac{\sigma^2}{1-\phi^2}} - \frac{n-1}{2}\log(2\pi) - \frac{n-1}{2}\log\left(\sigma^2\right) - \sum_{i=2}^n \frac{(x_i-\phi x_{i-1})^2}{2\sigma^2}$$
 treat X_1 as deterministic $= -\frac{n-1}{2}\log(2\pi) - \frac{n-1}{2}\log\left(\sigma^2\right) - \sum_{i=2}^n \frac{(x_i-\phi x_{i-1})^2}{2\sigma^2}$

Then, we have

$$\hat{\Psi}_{MLE} = \arg\max_{\Psi} \log L(\Psi)$$

For the coding work on this problem, check the attached jupyter notebook Time Series and Statistical Arbitrage HW2.ipynb

(b) Conditional MLE for MA(1) Process

Consider observations $\{X_1, X_2, \dots, X_n\}$ of a Gaussian MA(1) process. For parameter vector is $\Psi = (\theta_0, \theta_1, \sigma^2)$, we have

$$X_t = \theta_0 + \theta_1 Z_{t-1} + Z_t, \quad Z_t \sim IIDN(0, \sigma^2)$$

The probability distribution for X_t can be written as,

$$X_t \mid Z_{t-1} \sim \mathcal{N}\left(\theta_0 + \theta_1 z_{t-1}, \sigma^2\right) \Longrightarrow f\left(x_t \mid z_{t-1}\right) = \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left\{\frac{-\left(x_t - (\theta_0 + \theta_1 z_{t-1})\right)^2}{2\sigma^2}\right\}$$

Plug these in Equation (2), we obtain the Log-Likelihood function:

$$\log L(\Psi) = \log f(x_1 \mid \Psi) + \sum_{i=2}^n \log f(x_i \mid x_1, \dots, x_{i-1}, \Psi)$$
 treat X_1 as deterministic $= -\frac{n-1}{2} \log(2\pi) - \frac{n-1}{2} \log\left(\sigma^2\right) - \sum_{i=2}^n \frac{\left(x_t - \theta_0 - \theta_1 z_{t-1}\right)^2}{2\sigma^2}$

Maximize this function to estimate the parameters. (Much easier to implement)

Example: Suppose we have MA(1) process $X_t = 0.5Z_{t-1} + Z_t, Z_t \sim IIDN\left(0, 0.5^2\right)$. We simulate this process and pretend that we do not know the true values for ϕ and σ . Then, we use the conditional MLE to estimate the parameter vector $\Psi = (\theta_1, \sigma^2)$ By the above deduction, we get the log likelihood function is

$$\log L(\Psi) = -\frac{n-1}{2}\log(2\pi) - \frac{n-1}{2}\log(\sigma^2) - \sum_{i=2}^{n} \frac{(x_t - \theta_1 z_{t-1})^2}{2\sigma^2}$$

Then, we have

$$\hat{\Psi}_{MLE} = \arg\max_{\Psi} \log L(\Psi)$$

For the coding work on this problem, check the attached jupyter notebook Time Series and Statistical Arbitrage HW2.ipynb

3 Topic: MLE for AR and MA processes

Problem 3.1. For the time series model below:

$$X_t = 0.9X_{t-1} + Z_t$$
 Z_t is $N(0, 0.7^2)$

- (I) Simulate a path with 1000 data points.
- (II) Take the simulated path as the realized data and estimate the parameters using the MLE estimator.
- (III) Repeat steps (I) and (II) for 100 times and plot the distributions of the parameters you have estimated. What are the 95% confidence levels around your mean estimated parameters?
- (IV) Repeat (III) but use 5000 times instead of 100 times.

Use any programming language or algorithm you are comfortable with but we need to see the code.

Solution: Please check the attached jupyter notebook Time Series and Statistical Arbitrage HW2.ipynb

4 Application: Mean Squared Prediction and the Best Linear Predictor

Problem 4.1. An MA(1) process is given by

$$X_t = Z_t + \theta Z_{t-1}$$
 $\theta = 0.8$, $Z_t \sim WN(0, 1)$

Our observations indicates that $\{X_1 = 3.2020, X_2 = 1.5625\}$.

What is the best linear prediction and MSPE for X_3 ?

Solution: From Equation (1), we know that

$$P_{n}X_{n+h} := a_{0}^{*} + \mathbf{a}_{n}^{*^{\top}}\mathbf{x}_{n} \text{ where } \begin{cases} a_{0}^{*} &= \mu(1 - \mathbf{a}_{n}^{*^{\top}}\mathbb{1}) \\ \mathbf{a}_{n}^{*} &= \Gamma_{n}^{-1}\boldsymbol{\gamma}_{n}(h) \end{cases} \xrightarrow{n=2,h=1} P_{2}X_{3} := a_{0}^{*} + \mathbf{a}_{2}^{*^{\top}}\mathbf{x}_{2} \text{ where } \begin{cases} a_{0}^{*} &= \mu(1 - \mathbf{a}_{2}^{*^{\top}}\mathbb{1}) \\ \mathbf{a}_{2}^{*} &= \Gamma_{2}^{-1}\boldsymbol{\gamma}_{2}(1) \end{cases}$$

Since
$$\mathbb{E}[X_t] = 0 = \mu$$
 and $\gamma_X(h) = (1 + \theta^2)\gamma_Z(h) + \theta\gamma_Z(h-1) + \theta\gamma_Z(h+1) = \begin{cases} (1 + \theta^2)\sigma^2 & \text{if } h = 0 \\ \theta\sigma^2 & \text{if } h = \pm 1 \end{cases}$, we have otherwise

$$\Gamma_2 = \begin{bmatrix} \gamma_X(0) & \gamma_X(1) \\ \gamma_X(1) & \gamma_X(0) \end{bmatrix} = \begin{bmatrix} (1+\theta^2)\sigma^2 & \theta\sigma^2 \\ \theta\sigma^2 & (1+\theta^2)\sigma^2 \end{bmatrix}, \\ \gamma_2(1) = \begin{bmatrix} \gamma_X(1) \\ \gamma_X(2) \end{bmatrix} = \begin{bmatrix} \theta\sigma^2 \\ 0 \end{bmatrix} \\ \Longrightarrow \begin{cases} a_0^* & = 0 \\ \mathbf{a}_2^* & = \Gamma_2^{-1}\boldsymbol{\gamma}_2(1) = \frac{1}{1+\theta^2+\theta^4} \begin{bmatrix} \theta+\theta^3 \\ -\theta^2 \end{bmatrix} & \underline{\underline{\theta}=0.8} & \begin{bmatrix} \frac{820}{1281} \\ -\frac{400}{1281} \end{bmatrix} \end{cases}$$

Therefore, the Best Linear Predictor for X_3 is

$$\hat{X}_3 = P_n X_{n+h} := a_0^* + \mathbf{a}_n^{*^{\top}} \mathbf{x}_n = \frac{820}{1281} X_2 - \frac{400}{1281} X_1 = 1.5618$$

The corresponding MSPE is

$$MSPE = \mathbb{E}\left[\left(X_{n+h} - P_n X_{n+h}\right)^2\right] = \gamma_X(0) - \mathbf{a}_n^{*^{\top}} \boldsymbol{\gamma}_n(h) \xrightarrow{n=2,h=1} = \gamma_X(0) - \mathbf{a}_2^{*^{\top}} \boldsymbol{\gamma}_2(1) = 1.64 - \begin{bmatrix}\frac{820}{1281} & -\frac{400}{1281}\end{bmatrix} \begin{bmatrix}0.8\\0\end{bmatrix} = \frac{656}{1281}$$

Hence, the Best Linear Predictor for X_3 is $\hat{X}_3 = 1.5618$ and the corresponding $MSPE = \frac{656}{1281}$.