Project 1

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Problem 1

Part A

For Part A, the following statistics were calculated:

Mean: 0.050198Variance: 0.010322Skewness: 0.120626Kurtosis: 0.230070

Part B

Part B compared a Normal Distribution with a T-Distribution. The observed kurtosis was higher than 0, suggesting heavier tails than a normal. Based on this, the T-Distribution would be preferable.

Part C

Both distributions were fitted using MLE. The more negative AIC and BIC values for the Normal Distribution indicated a better fit, relative to the initial kurtosis result. Therefore, a Normal Distribution was chosen as the better fit.

Problem 2

Part A: Pairwise Covariance Matrix

A pairwise approach was used to compute covariances between columns, ignoring rows lacking data for each pair. The matrix obtained was:

| 1.47048437 | 1.45421424 | 0.87726904 | 1.90322645 | 1.44436105 |
|------------|------------|------------|------------|------------|
| 1.45421424 | 1.25207795 | 0.53954816 | 1.62191837 | 1.23787697 |
| 0.87726904 | 0.53954816 | 1.272425 | 1.17195897 | 1.091912 |
| 1.90322645 | 1.62191837 | 1.17195897 | 1.81446921 | 1.58972858 |
| 1.44436105 | 1.23787697 | 1.091912 | 1.58972858 | 1.39618646 |

Part B: Positive Semi-Definiteness Check

The eigenvalues were:

Eigenvalue

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| Eigenvalue |
|-------------|
| 6.78670573 |
| 0.83443367 |
| -0.31024286 |
| 0.02797828 |
| -0.13323183 |

Two negative eigenvalues were found, indicating the matrix is not positive semi-definite.

Part C: Nearest PSD Matrix

Two methods were used: Rebenato and Jackel, and Higham's method.

Rebenato and Jackel

This approach sets all negative eigenvalues to zero after eigen-decomposition and reconstructs the matrix:

| 1.61513295 | 1.44196041 | 0.89714421 | 1.78042572 | 1.43379434 |
|------------|------------|------------|------------|------------|
| 1.44196041 | 1.34696791 | 0.58508635 | 1.55455193 | 1.21140918 |
| 0.89714421 | 0.58508635 | 1.29891578 | 1.11595578 | 1.07669234 |
| 1.78042572 | 1.55455193 | 1.11595578 | 1.98316488 | 1.62137332 |
| 1.43379434 | 1.21140918 | 1.07669234 | 1.62137332 | 1.40493616 |

Higham's Method

- 1. With matrix A, set $Y_0 = A^{**}$, $\Delta S_0 = 0$.
- 2. $R_k = Y_{k-1} \Delta S_{k-1}$.
- 3. Clamp negative eigenvalues of R_k to zero.
- 4. Update ΔS.
- 5. Repeat until the difference $||Y_k Y_{k-1}||$ is below a tolerance or the maximum number of iterations is reached.

| 1.61513295 | 1.44196041 | 0.89714421 | 1.78042572 | 1.43379434 |
|------------|------------|------------|------------|------------|
| 1.44196041 | 1.34696791 | 0.58508635 | 1.55455193 | 1.21140918 |
| 0.89714421 | 0.58508635 | 1.29891578 | 1.11595578 | 1.07669234 |
| 1.78042572 | 1.55455193 | 1.11595578 | 1.98316488 | 1.62137332 |
| 1.43379434 | 1.21140918 | 1.07669234 | 1.62137332 | 1.40493616 |

Part D: Overlapping Data Covariance

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Excluding rows with missing data for any column produced the following covariance matrix (using same cov matrix methodology from part A):

| 0.41860366 | 0.39405407 | 0.42445735 | 0.41638241 | 0.43428682 |
|------------|------------|------------|------------|------------|
| 0.39405407 | 0.39678563 | 0.40934344 | 0.3984012 | 0.42263077 |
| 0.42445735 | 0.40934344 | 0.4413601 | 0.42844141 | 0.44895733 |
| 0.41638241 | 0.3984012 | 0.42844141 | 0.43727358 | 0.44016735 |
| 0.43428682 | 0.42263077 | 0.44895733 | 0.44016735 | 0.46627249 |

Part E: Comparison of the Matrices from (C) and (D)

The matrices in (C) use maximum available pairs, which can create inconsistencies leading to negative eigenvalues. The overlapping-data matrix in (D) is PSD by construction but may discard many rows. The off-diagonal values in the near-PSD matrices can differ significantly from the overlapping-data approach because of sample size differences.

Problem 3

Part A: Fitting a Multivariate Normal

After loading the dataset, the sample mean vector and covariance matrix were computed for X_1 and X_2 . The mean is approximately [0.0460, 0.0999], and the covariance matrix is:

| 0.0101622 | 0.00492354 |
|------------|------------|
| 0.00492354 | 0.02028441 |

Part B: Distribution of X_2 Given $X_1 = 0.6$

Method 1 (Conditional Formula) Using the standard bivariate normal conditional formula gave:

- Conditional Mean ~ 0.368325
- Conditional Variance ~ 0.017899

Method 2 (Linear Regression) A simple linear regression of X_2 on X_1 yielded:

- Conditional Mean ~ 0.368325
- Residual MSE ~ 0.017917

Both methods align well.

Part C: Cholesky-Based Simulation

A large sample (X_1, X_2) was generated from $N(\mu, \Sigma)$. After filtering for X_1 close to 0.6:

- Simulated Conditional Mean ~ 0.368784
- Simulated Conditional Variance ~ 0.014673

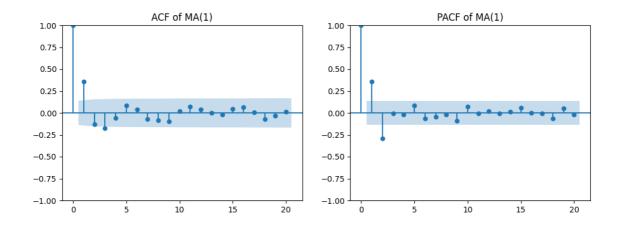
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These are consistent with the analytical approach.

Problem 4

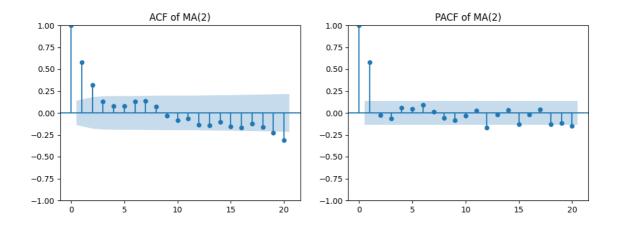
Part A: Simulating MA(1), MA(2), and MA(3)

1. MA(1)



- ACF: Notice a significant spike at lag 1, then it drops near zero for higher lags.
- PACF: The first two lags are significant. Lag 17 is also significant.

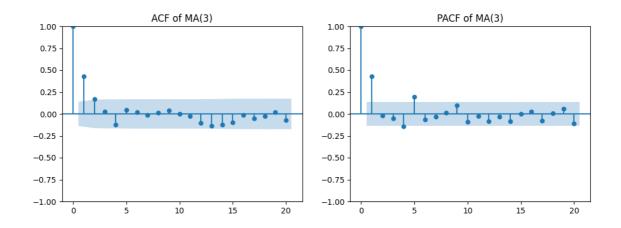
2. MA(2)



- ACF: Typically shows two significant spikes (lags 1 and 2) before going near zero for larger lags.
- **PACF**: Although MA(1) doesn't exhibit gradual tapering, MA(2) PCF shows an even more aggressive drop after the first lag.

3. **MA(3)**

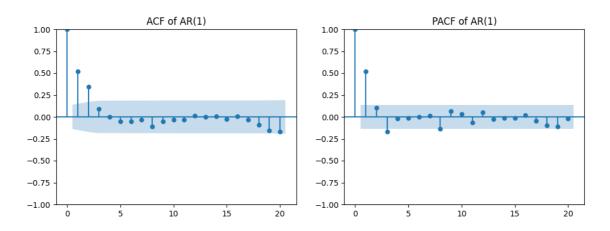
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- ACF: Three significant autocorrelation spikes (lags 1, 2, and 3), dropping off afterward.
- PACF: Similar to MA(2) PCF, gradual tapering isn't seen.

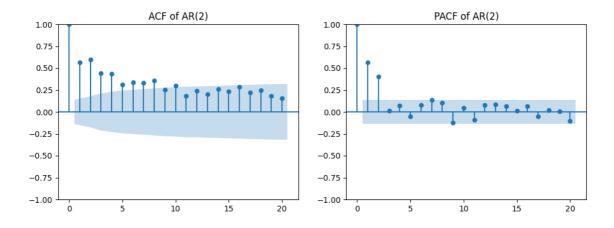
Part B: Simulating AR(1), AR(2), and AR(3)

1. **AR(1)**



- ACF: It tails off exponentially, but it cuts off after lag 2.
- **PACF**: Shows a clear significant spike at lag 1, then becomes very small afterward.

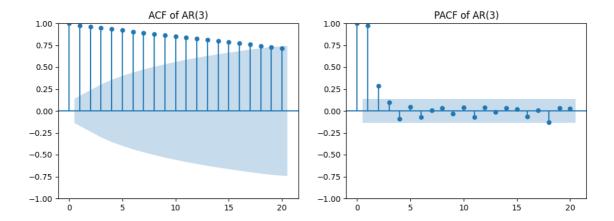
2. AR(2)



- **ACF**: It decays as expected for an AR model.
- PACF: Significant spikes at lags 1 and 2, then mostly zero for higher lags.

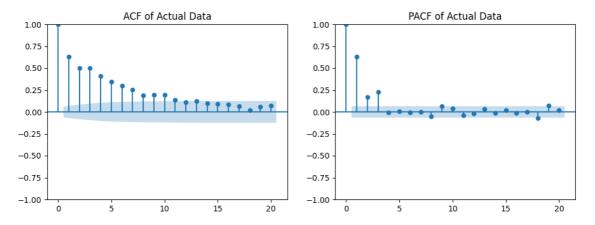
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3. AR(3)



- **ACF**: It decays, but it decays at an even slower rate than seen in AR(20).
- **PACF**: Three significant partial autocorrelation spikes, followed by lower values afterward. This confirms AR(3) structure.

Part C: Examining problem4.csv



The ACF gradually declines, and the PACF has three prominent lags. This suggests an AR(3) model.

Part D

Model fit comparison using AIC and AICc:

```
(1, 0, 0) => (AIC=-1669.089, AICc=-1669.065)

(2, 0, 0) => (AIC=-1696.092, AICc=-1696.051)

(3, 0, 0) => (AIC=-1746.282, AICc=-1746.221)

(0, 0, 1) => (AIC=-1508.927, AICc=-1508.903)

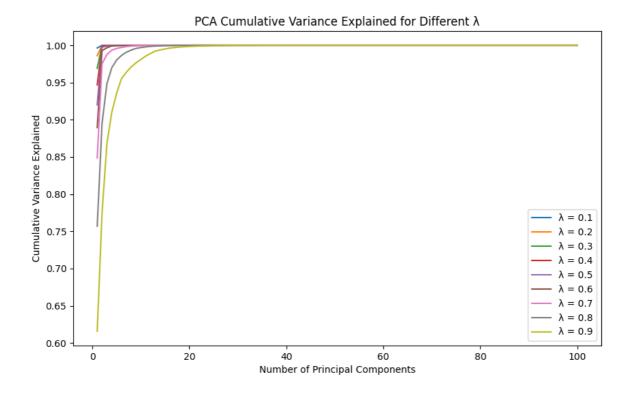
(0, 0, 2) => (AIC=-1559.251, AICc=-1559.211)

(0, 0, 3) => (AIC=-1645.133, AICc=-1645.073)
```

The best model by AIC is AR(3).

Problem 5

Part B



Part C

Larger lambda values decay more slowly, so older observations remain significant and the covariance matrix is more stable. Smaller lambda values make recent data more influential, producing a sharper drop in the leading principal components.

Problem 6

Part A

A Cholesky Root simulation was performed with a near-psd matrix for stability. The matrix was generated using the near-psd method of Rebenato and Jackel.

Part C

The Frobenius norm relative to the original is 0.000006 for Cholesky-based and 0.000022 for PCA-based, indicating the Cholesky-based matrix is closer.

Part D

The Cholesky-based matrix preserves the entire covariance structure, leading to a gradual cumulative variance. The PCA-based method explains 66% by the first component, 84% by the seventh, and 100% by the 29th, reflecting fewer non-zero eigenvalues and a lower rank.

Part F: Tradeoffs

The Cholesky approach is quicker and retains the full covariance. PCA offers dimension reduction and noise filtering, which can be beneficial in high-dimensional settings or when dropping minor components is preferred.