ADSA Data Analytics Project — Final Report

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1. Project overview and objective

Goal. Build a reproducible ML pipeline that ingests movie (or given) metadata, performs EDA, cleans and transforms the data, trains and validates models, and exposes a prediction pipeline. Deliverables: working code (src/), documented GitHub repo, reproducible environment, and a short report containing results and next steps.

Data. Describe dataset(s) used (columns, number of rows, source). If multiple files (CSV/BigQuery), list them here.

Success criteria. e.g., target metric (RMSE, MAE, R2, Accuracy), baseline performance, and production readiness requirements (unit tests, logging, packaging).

2. Step 1 — GitHub & code setup (expanded)

2.1 Environment creation (conda + venv)

Use an isolated environment to ensure reproducibility. Create a conda environment in the project folder and activate it.

Typical commands to run:

```
    conda create -p ./venv python==3.8 -y
    conda activate ./venv
    Install dependencies from requirements.txt after creating it: pip install -r requirements.txt
```

Note: Use explicit pinned versions in requirements.txt for reproducibility.

2.2 Git initialization & remote

Initialize a repository, create README and .gitignore, add files, commit, set the main branch and add a remote.

Typical git commands to run:

```
    git init
    git add README.md .gitignore
    git commit -m "first commit"
    git branch -M main
    git remote add origin https://github.com/Manoj-Sh-AI/ADSA_Data_Analytics-1_ml_pipeline.git
    git push -u origin main
```

2.3 Project packaging (setup.py)

If you plan to install the project as a package during development, create a src/ package and a minimal setup.py. Use pip install -e . to install locally during development.

2.4 Requirements

Keep requirements.txt updated and pin versions. Example packages to include: pandas, numpy, scikit-learn, xgboost, catboost, joblib, pyyaml, pytest.

3. Step 2 — Error handling, logging & project structure

3.1 Project structure (high-level)

(See Appendix A for full skeletal tree.)

Key folders and responsibilities:

- src/components/: modular units (data_ingestion, data_transformation, model_trainer)
- src/pipeline/: orchestrates end-to-end flow (train_pipeline, predict_pipeline)
- src/config/: YAML/JSON with paths, hyperparameters, and constants
- src/logs/: runtime logs
- tests/: unit & integration tests

3.2 Logging

Use Python logging configured to write to both console and a file in a logs/ folder. Provide a helper (e.g., get_logger) that sets up formatters, stream handler and file handler. Call the logger at the top of each module to standardize messages and timestamps.

3.3 Exception handling

Create a custom exception class to attach context and preserve tracebacks. Wrap high-level pipeline functions with try/except blocks, log the exception with stack trace (using logger.exception), and re-raise or handle with meaningful messages.

4. Step 3 — Data Ingestion

4.1 Goals

- Read raw data from CSV / BigQuery / local storage
- Validate schema (expected columns)
- Save raw snapshot (versioned) and a cleaned copy for transformations

4.2 Data ingestion API (recommended behaviour)

Design data ingestion to:

- Accept configuration (paths for raw, train and test)
- Read the raw dataset
- Perform a deterministic train/test split (e.g., 80/20, with a fixed seed)
- Persist the train and test sets to data/processed/
- · Log operations and raise meaningful exceptions on failure

4.3 Validation checks

- Column presence
- Missing value proportions per column
- Duplicate rows
- Data types and ranges (e.g., dates, numeric ranges)

5. Step 4 — Exploratory Data Analysis (EDA)

5.1 Objectives

- Understand distributions, missingness, correlation with target
- · Detect outliers and data quality issues
- Identify candidate feature transformations

5.2 Recommended EDA checklist

- Summary statistics (describe and info)
- Missing value heatmap and percent-missing table
- Univariate plots for numeric features (histograms / boxplots)
- Countplots for categorical variables
- Correlation matrix / heatmap for numeric relationships
- Target vs feature plots (scatter, violin, bar) for top predictors
- Time-series checks if data is temporal

5.3 Tools

Pandas and Matplotlib (or Seaborn for convenience) are recommended. Conduct EDA interactively in Jupyter notebooks and save relevant visuals to reports/figures/.

6. Step 5 — Data Transformation & Feature Engineering

6.1 Goals

- · Convert raw features into model-ready features
- · Handle missing data, encode categoricals, scale numeric features
- Persist transformers (scalers, encoders) with joblib for inference

6.2 Typical steps

- 1. Imputation: numeric (mean/median/KNN), categorical (mode or new category "Missing")
- 2. **Categorical encoding**: One-Hot for low-cardinality, Target/Ordinal or Count encoding for high-cardinality
- 3. Scaling: Standard scaling for models sensitive to scale, MinMax for neural networks
- 4. Feature creation: interaction terms, datetime decompositions, aggregated features
- 5. Dimensionality reduction (if needed): PCA, feature selection via tree-based importance

6.3 Implementation notes

- Build a preprocessing pipeline object that encapsulates imputation, encoding and scaling and fit it on the training data only.
- Persist the fitted preprocessing object to <u>artifacts</u>/ so it can be reused by the prediction pipeline.

7. Step 6 — Model Training (Why I chose these models and how I achieved the accuracy)

7.1 Modeling strategy

- Try multiple models: Linear Regression (baseline), RandomForest, XGBoost, CatBoost, and a simple neural net if needed.
- Use cross-validation (KFold, or TimeSeriesSplit if temporal) and compare with metrics (RMSE, MAE, R2). Keep a validation set for final selection.

7.2 Why these models?

- Linear Regression: interpretable baseline, fast.
- RandomForest: strong baseline for tabular data, robust to outliers and feature scaling.
- **XGBoost/CatBoost**: state-of-the-art for many tabular problems; often give best performance with moderate tuning.

7.3 Hyperparameter tuning

Use randomized or grid search for hyperparameter tuning; for faster, use Bayesian optimizers (Optuna). Keep $\begin{bmatrix} n_j obs=-1 \end{bmatrix}$ where possible and use sensible search spaces.

7.4 Evaluation

Report train/validation/test metrics and show learning curves. Record metrics in a reports/ file and compare several candidate models in a table.

7.5 Persisting the model

Persist the trained model artifact to <code>artifacts/</code> (e.g., <code>model.joblib</code>) for later inference.

8. Step 7 — Prediction / Inference pipeline

8.1 Design

- Provide a single entry point (predict pipeline) that accepts raw input (CSV or JSON), applies the saved preprocessing object, loads the model artifact, and outputs predictions (CSV/JSON).
- Validate input schema and produce outputs with predictable column names (e.g., prediction).

8.2 Inference behaviour

- Load preprocessor and model artifacts from artifacts/.
- Transform the raw input using the preprocessor and call model.predict() .
- Return or persist predictions alongside input identifiers.

8.3 Deployment considerations

- For a demo: expose via a lightweight API (Flask/FastAPI).
- For production: containerize, add health checks, secure secrets, and add monitoring.

9. Reproducibility & How to run the pipeline

9.1 Quick start (development)

Typical steps to run locally:

- 1. Clone the repo.
- 2. Create the environment and install requirements.
- 3. conda create -p ./venv python==3.8 -y
- 4. conda activate ./venv
- 5. pip install -r requirements.txt
- 6. Run the training pipeline (script entry point) and the prediction script with input and output paths. Ensure config/ contains correct paths and seeds.

10. Results summary & discussion

Include here a short table with final metrics for the chosen model and the baseline. Also include: feature importances, top 5 features, and saved visuals (learning curves, residual plots) placed in reports/figures/.

Example columns to report: Model, Dataset, RMSE, MAE, R2.

11. Conclusion, limitations & future work

- Conclusion: Which model you selected and why, summary of results.
- Limitations: data quantity/quality, label noise, potential leakage, or limited features.
- **Future work:** more feature engineering, ensembling, better cross-validation, model explainability (SHAP), productionizing with CI/CD.

12. File Structure

Appendix A — Skeletal file structure

```
ADSA_Data_Analytics-1_ml_pipeline/

├ .gitignore
├ README.md
├ requirements.txt
⊢ setup.py
├ config/
   └ config.yaml
 - data/
   ⊢ raw/
   ⊢ processed/
   └ external/
 - src/
   ⊢ __init__.py
   ⊢ logger.py
   ⊢ exception.py
   ⊢ utils.py

    ⊢ components/

├ __init__.py

    ⊢ data_ingestion.py

      └ model_trainer.py
   ⊢ pipeline/
      ├ __init__.py

    ⊢ train_pipeline.py

      └ predict_pipeline.py
   └ config/
      └ params.yaml
├ artifacts/
   ⊢ preprocessor.joblib
   └ model.joblib
⊢ reports/
```

```
| ├─ figures/
| └─ report.pdf
└─ tests/
└─ test_data_ingestion.py
```

Appendix B — Sample scripts & snippets

- train_pipeline should: load config → run data_ingestion → run data_transformation → run model_trainer → save artifacts and metrics.
- predict_pipeline should: parse args → validate input → load artifacts → run predict → save results.