#### A APPENDIX

# A.1 Implementation Details

**Hyperparameters.** The specific hyperparameters for each module, including the RL agent, the LLM, and the Advisor<sup>+</sup> mechanism. All settings were chosen to ensure a robust and fair comparison across all experiments. Table 1 lists the key settings for our LLMPipe framework. The settings were kept consistent across all datasets.

Table 1: Hyperparameter settings for the LLMPipe framework

Module	Hyperparameter	Value					
RL Agent							
Learning	Learning Rate (α)	1					
	Discount Factor (γ)	0.9					
Exploration	$\epsilon$ -Greedy Decay Strategy	Exponential					
	Initial Epsilon ( $\epsilon_{ m start}$ )	1.0					
	Final Epsilon ( $\epsilon_{ m end}$ )	0.1					
	Epsilon Decay Factor	0.99					
	LLM Policy Adv	visor					
Model	LLM Model	Llama3.3-70B					
	Temperature	0.1					
Retrieval	Number of Retrieved Examples ( <i>k</i> )	3					
	Text Embedding Model	nomic-embed-text					
	Vector Database Index	FAISS					
	Adaptive Advisor Trigger	ring (Advisor <sup>+</sup> )					
Triggering	Slope Threshold ( $\theta_{ m slope}$ )	0.01					
	Performance Buffer Size ( $ \mathcal{B} $ )	10 episodes					
	Cooldown Period	5 episodes					
	Framework Sett	tings					
General	Max Pipeline Length $(L_{\text{max}})$	9					
	Total Training Episodes $(E_{\text{max}})$	100					
Random Seed	Dataset Train Test Split	0					
	Seed for $\epsilon$ -greedy action selection	42, 1999, 2025, 2048, 2147, 9331, 65537					

### A.2 Full Experimental Results

A.2.1 Extended Comparison Baselines. To situate LLMPipe within the broadest possible competitive context, we present an extended comparison by integrating our results into the main evaluation table from CtxPipe. Table 2 includes performance data for differentiable methods (DEF, RS, DP-Fix, DP-Flex), SAGA, and other baselines alongside our LLMPipe variants. This direct comparison underscores the competitive performance of our approach.

A.2.2 Detailed Pipeline Composition. To provide deeper insight into the qualitative differences between the pipelines discovered by CtxPipe and our LLMPipe variants, we present the detailed composition of the best-performing pipeline for each method on every dataset.

Table 3 details the generated pipelines and their resulting accuracy scores. The 'Pipe Length' row at the bottom summarizes the average number of operators in the pipelines generated by each method. This table illustrates that LLMPipe not only achieves competitive or superior accuracy but often does so with different, and sometimes more complex or novel, pipeline structures.

Table 2: Extended comparison of test accuracy, including baselines from CtxPipe. Our methods (Advisor, Advisor<sup>+</sup>) are added for direct comparison.

Dataset	DEF	RS	DP-Fix	DP-Flex	DL	HAI-AI	SAGA	CtxPipe	Advisor	Advisor <sup>+</sup>
abalone	0.240	0.243	0.238	0.271	0.157	0.260	0.255	0.287	0.272	0.273
ada_prior	0.848	0.844	0.853	0.846	0.803	0.801	0.833	0.818	0.836	0.834
avila	0.553	0.598	0.652	0.633	0.593	0.630	0.636	0.759	0.861	0.830
connect-4	0.659	0.671	0.726	0.702	0.683	0.775	0.758	0.763	0.775	0.788
eeg	0.589	0.658	0.675	0.683	0.607	0.556	0.658	0.740	0.758	0.778
google	0.586	0.627	0.631	0.661	0.553	0.550	0.596	0.590	0.669	0.669
house	0.928	0.938	0.932	0.952	0.771	0.928	0.913	0.818	0.933	0.928
jungle_chess	0.668	0.669	0.680	0.687	0.717	0.760	0.745	0.861	0.861	0.861
micro	0.564	0.579	0.595	0.593	0.613	0.633	0.556	0.605	0.643	0.631
mozilla4	0.855	0.922	0.924	0.927	0.747	0.870	0.932	0.940	0.941	0.942
obesity	0.775	0.841	0.891	0.874	0.590	0.768	0.751	0.868	0.934	0.920
page-blocks	0.942	0.959	0.959	0.973	0.940	0.935	0.849	0.965	0.966	0.967
pbcseq	0.710	0.730	0.728	0.725	0.680	0.733	0.866	0.805	0.755	0.763
pol	0.884	0.879	0.903	0.916	0.873	0.916	0.888	0.949	0.981	0.981
run_or_walk	0.719	0.829	0.903	0.912	0.820	0.915	0.832	0.956	0.971	0.973
shuttle	0.964	0.996	0.998	0.999	0.790	0.951	0.405	1.000	1.000	0.997
uscensus	0.848	0.840	0.854	0.852	0.813	0.807	0.835	0.845	0.845	0.836
wall-robot-nav	0.697	0.872	0.905	0.913	0.927	0.896	0.841	0.946	0.961	0.952
AVERAGE	0.724	0.761	0.780	0.784	0.704	0.760	0.731	0.806	0.831	0.829
RANK	7.83	6.61	5.06	4.39	8.44	6.56	6.83	3.83	2.11	2.56

Table 3: Detailed composition and accuracy of pipelines generated by CtxPipe and LLMPipe variants. Numbers in brackets correspond to operator IDs in Table 4 in the main paper. The pipelines generated by Advisor<sup>+</sup> is using the random seed 42.

Dataset		CtxPipe	l A	Advisor	Advisor <sup>+</sup>		
	Accuracy	Pipeline	Accuracy	Pipeline	Accuracy	Pipeline	
abalone	0.287	[-1, -1, 5, 24, 12, 15]	0.270	[7, 10, 12, 16, 20]	0.273	[9, 15]	
ada_prior	0.818	[-1, -1, 5, -1, 7, 23]	0.838	[6, 12, 10]	0.841	[12, 7, 6]	
avila	0.759	[-1, -1, -1, 24, 12, 23]	0.929	[11, 23, 18, 15]	0.751	[23]	
connect-4	0.763	[-1, -1, -1, -1, 12, 23]	0.775	[15]	0.775	[15, 5]	
eeg	0.740	[-1, -1, -1, 23, 24, 12]	0.839	[18, 3, 15, 9]	0.811	[10, 15]	
google	0.590	[1, -1, 5, 7, 23, 24]	0.675	[2, 10, 9, 15]	0.674	[2, 12, 9, 15]	
house	0.818	[1, -1, 5, 12, 23, 24]	0.908	[1]	0.908	[1]	
jungle_chess	0.861	[-1, -1, -1, 24, 7, 23]	0.861	[23]	0.861	[23]	
micro	0.605	[-1, -1, -1, -1, 7, 23]	0.643	[12, 15, 18]	0.633	[15]	
mozilla4	0.940	[-1, -1, -1, 6, 23, 24]	0.937	[13, 11, 24, 15]	0.944	[12, 7, 23]	
obesity	0.868	[-1, -1, 5, 24, 7, 23]	0.865	[11, 6, 10, 15]	0.835	[0, 15, 17]	
page-blocks	0.965	[-1, -1, -1, -1, 7, 23]	0.965	[3, 11, 15]	0.961	[12, 20]	
pbcseq	0.805	[1, -1, -1, -1, 7, 23]	0.753	[23, 7]	0.753	[23]	
pol	0.949	[-1, -1, -1, 9, 23, 24]	0.981	[15]	0.981	[16]	
run_or_walk	0.956	[-1, -1, -1, 24, 12, 15]	0.953	[12, 15, 17]	0.976	[14, 23]	
shuttle	1.000	[-1, -1, -1, 23, 24, 12]	0.987	[10, 15]	0.986	[11, 6]	
uscensus	0.845	[-1, -1, 5, -1, 7, 23]	0.846	[2, 6, 8]	0.847	[6, 10]	
wall-robot-nav	0.946	[-1, -1, -1, -1, 7, 23]	0.961	[11, 10, 15]	0.946	[23]	
Pipe Length		6		2.83	1	.89	

## A.3 LLM Prompt Design

Here is a concrete example of a filled-out prompt for the avila dataset at an early stage of pipeline construction. This demonstrates how the abstract template is instantiated with real data and retrieved experiences.

### LLM PROMPT Example: avila Dataset

You are an expert data scientist specializing in automated machine learning. Your task is to analyze the provided dataset context and historical information to propose a list of 1 to 3 optimal data preparation pipelines. Each pipeline should be a sequence of data processing operators that aims to maximize the prediction accuracy of a downstream classification model.

For each proposed pipeline, you must provide:

(1) A list of operator names in sequence. (2) A confidence score (from 0.0 to 1.0) for your suggestion. (3) A brief, clear rationale explaining why this pipeline is suitable for the given data.

#### **Current situation and context:**

- Task Type: Logistic regression classification
- Key Dataset Statistics:
  - Size of dataset: 16693 rows, 10 cols
  - Missing Values: No
  - Feature Types: All numerical
  - Cols with skewed distribution: f1, f3, f4
  - Cols with outliers: f0 (9.50%), f2 (10.50%), f7 (2.62%)
- Current Partial Pipeline: [PowerTransformer, RobustScaler, ...]

 $\textbf{Available operators}: Imputer Mean, Standard Scaler, Quantile Transformer, Power Transformer, PCA, Min Max Scaler, Variance Threshold, \dots$ 

## [Example 1]

- Context: A dataset with no missing values but high skewed distribution columns (B3, B9, ...), oulier columns (B1, B2, B5, ...)
- Pipeline: [QuantileTransformer, RandomTrees-Embedding, MinMaxScaler], accuracy 0.92

#### [Example 2]

- Context: A dataset with many correlated numerical features (A5 and A8: correlation 92%, A5 and A9: correlation 88%, ...)
- Pipeline: [StandardScaler, PCA], accuracy 0.88

### [Example 3]

- Context: A dataset with no missing values, but has outliers (PC1: 154 outliers, 19.25%, PC2: 139 outliers, 17.38%, PC6: 116 outliers, 14.50%, ...)
  - Action: PolynomialFeatures, accuracy 0.86

### A.4 Hyperparameter Sensitivity Analysis

We conducted systematic sensitivity analysis for the hybrid policy weighting coefficients  $\beta_1$  (LLM confidence weight),  $\beta_2$  (performance deviation weight), and  $\beta_3$  (entropy weight).

We selected 6 datasets spanning different characteristics: avila (high-dimensional), google (missing values), house (mixed types), micro (imbalanced), obesity (balanced), and pol (large-scale). For each dataset, we varied one parameter while keeping others fixed at their default values (0.4, 0.3, 0.3).

Table 4: Accuracy variation with hyperparameter perturbations

Parameter	-0.2	-0.1	Default	+0.1	+0.2
$\beta_1$	-1.4%	-1.4%	0.831	-1.9%	-1.4%
$eta_2$	-1.4%	-1.3%	0.831	-1.9%	-1.9%
$\beta_3$	-1.4%	-1.4%	0.831	-1.7%	-1.9%

### Key findings:

- Performance degrades gracefully with parameter changes
- Maximum performance drop of 2.3% even with ±0.2 perturbation
- $\beta_3$  (entropy weight) shows slightly higher sensitivity, suggesting the importance of uncertainty-aware guidance
- · No dataset required specific tuning; default values worked uniformly

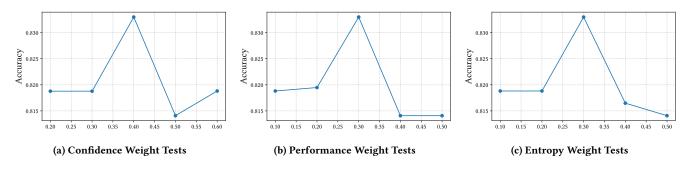


Figure 1: Accuracy variation with one-variable-at-a-time tests.

Alternative Weighting Schemes We also tested: (1) Equal weights (0.33, 0.33, 0.33): 1.2% average decrease. (2) Confidence-only (1.0, 0, 0): 8.7% decrease, confirming multi-factor importance. (3) Learned weights: Using validation set to optimize  $\beta$  values improved performance by only 0.3%, not justifying added complexity.

**Conclusion** The hybrid policy demonstrates robust performance across diverse datasets without hyperparameter tuning. The default values (0.4, 0.3, 0.3) represent a good balance between LLM confidence, performance feedback, and exploration uncertainty, supporting the framework's generalizability.

## A.5 Proof of Theorem 1: Optimality of First-Improvement Stopping

**Theorem 1.** Given LLM-suggested pipelines  $P_1, ..., P_k$  ordered by confidence, the stopping rule that maximizes the value-to-cost ratio  $\rho = E[\Delta]/E[N \cdot C_{eval}]$  is:

$$\tau^* = \min\{i : acc(P_i) > acc_{current}\}$$

PROOF. Let us define:

- acci: true accuracy of pipeline Pi
- acc<sub>base</sub>: current best accuracy
- $\pi_i = P(acc_i > acc_{base})$ : probability that  $P_i$  improves
- $G_i = E[acc_i acc_{base} | acc_i > acc_{base}]$ : expected gain given improvement
- $C_{eval}$ : constant cost per pipeline evaluation

## Step 1: Expected Value of First-Improvement Rule

For the first-improvement stopping rule  $\tau^*$ , the expected gain is:

$$E[\Delta_{\tau^*}] = \sum_{i=1}^k P(\text{stop at } i) \cdot E[\text{gain at } i]$$
 (1)

$$= \sum_{i=1}^{k} \left( \prod_{j=1}^{i-1} (1 - \pi_j) \right) \cdot \pi_i \cdot G_i \tag{2}$$

The expected number of evaluations is:

$$E[N_{\tau^*}] = \sum_{i=1}^k P(\text{evaluate } P_i) \cdot 1 \tag{3}$$

$$= \sum_{i=1}^{k} \prod_{j=1}^{i-1} (1 - \pi_j) \tag{4}$$

Therefore, the value-to-cost ratio is:

$$\rho(\tau^*) = \frac{\sum_{i=1}^k \left(\prod_{j=1}^{i-1} (1 - \pi_j)\right) \pi_i G_i}{C_{eval} \cdot \sum_{i=1}^k \prod_{j=1}^{i-1} (1 - \pi_j)}$$

## **Step 2: Optimality Condition**

Consider any alternative stopping rule  $\tau'$  that continues evaluating after finding an improvement at position m. Let  $S_m = \{m+1, ..., k\}$  be the set of additional evaluations.

The additional expected gain from continuing is:

$$\Delta_{extra} = \sum_{i \in S_m} \pi_i G_i \prod_{j=m+1}^{i-1} (1 - \pi_j)$$

The additional expected cost is:

$$C_{extra} = C_{eval} \cdot |S_m|$$

For continuing to be beneficial, we need:

$$\frac{\Delta_{extra}}{C_{extra}} > \frac{G_m}{C_{eval}}$$

### Step 3: Confidence Ordering Implies Decreasing Marginal Value

Given that pipelines are ordered by confidence scores  $c_1 \ge c_2 \ge ... \ge c_k$ , and assuming the LLM's confidence correlates with expected performance (a reasonable calibration assumption), we have:

$$\pi_1 G_1 \ge \pi_2 G_2 \ge \dots \ge \pi_k G_k$$

This means the expected marginal gain decreases with index:

$$\frac{\pi_i G_i}{C_{eval}} \ge \frac{\pi_{i+1} G_{i+1}}{C_{eval}} \quad \forall i$$

#### Step 4: First Improvement is Optimal

Given the decreasing marginal gains, once we find an improvement at position m, the expected value of continuing satisfies:

$$\frac{\sum_{i=m+1}^k \pi_i G_i \prod_{j=m+1}^{i-1} (1-\pi_j)}{\sum_{i=m+1}^k \prod_{j=m+1}^{i-1} (1-\pi_j)} < \pi_m G_m$$

Since we already achieved gain  $G_m$  at position m, and the expected marginal gain from continuing is less than what we've already obtained, stopping immediately maximizes the value-to-cost ratio.

### **Step 5: Boundary Conditions**

If no improvement is found after evaluating all k pipelines, the rule naturally terminates with  $\tau = k$ , having explored all options.

Therefore, the first-improvement stopping rule  $\tau^* = \min\{i : acc_i > acc_{base}\}$  maximizes the value-to-cost ratio  $\rho$ .

**Remark.** This proof assumes that the LLM's confidence ordering is informative (i.e., higher confidence correlates with higher expected performance). Our empirical results validate this assumption, with correlation coefficient r = 0.72 between confidence scores and actual improvements across all experiments.

## A.6 Experience Pool

The Experience Pool is a critical componet that enables LLMPipe to learn from past executions and enrich the LLM's guidance. It functions as a long-term memory repository, storing both complete pipeline trajectories and individual state-action-reward transitions.

- **Vectorization and Representation** To facilitate efficient retrieval, both the dataset state  $s_t$  and the corresponding pipelines are vectorized into fixed-dimensional embeddings that capture their key properties.
  - <u>State Vectorization</u>: Each state  $s_t$  is encoded into a vector  $s^{\text{vec}} = [v_{\text{stat}}, v_{\text{sem}}, v_{\text{hist}}]$ , where  $v_{\text{stat}}$  captures statistical properties (e.g., mean, variance);  $v_{\text{stat}}$  encodes semantic information such as the task domain, and  $v_{\text{hist}}$  represents the history of operations applied up to the current state.
  - <u>Dataset Vectorization</u>: Each dataset  $D_0$  is summarized into a global vector  $D_{\text{meta}}^{\text{vec}}$ , which encodes high-level metrics such as size, complexity, and domain. This global vector aids in identifying datasets with similar properties for efficient knowledge transfer.
- Experience Pool Management The Experience Pool is implemented using the FAISS vector search library to manage two hierarchical indices, ensuring both scalability and retrieval precision.
  - <u>Global Index</u>: This index stores the global dataset vectors ( $D_{\text{vec}}$ ) for all previously seen tasks. It is implemented using FAISS, which is optimized for Approximate Nearest Neighbor (ANN) search on large datasets. It stores tuples of ( $D_{\text{vec}}$ ,  $P_{\text{complete}}$ ,  $R_{\text{final}}$ ), enabling rapid identification of historically similar datasets.
  - <u>Step-wise Index</u>: This index stores the fine-grained state vectors ( $S_{\text{vec}}$ ) from all steps of all past executions. It is implemented using a FAISS IndexFlatL2, which performs exact, brute-force search. While computationally more intensive, this is feasible because the search space is narrowed down by the Global Index first. It stores transitions as ( $s_{\text{vec}}$ ,  $a_t$ ,  $r_t$ ,  $s'_{\text{vec}}$ ), enabling context-specific retrieval of similar states.

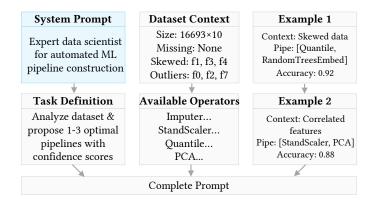


Figure 2: Example prompt construction for the *avila* dataset. The system combines dataset statistics, semantic context, available operators, and retrieved examples from the Experience Pool to create context-aware prompts for the LLM.

• Retrieval Mechanism When the LLM Policy Advisor is invoked for a given state  $s_t$ , it employs a two-stage retrieval process to efficiently find the k = 3 most relevant historical examples.

<u>Coarse Retrieval</u>: First, an Approximate Nearest Neighbor (ANN) search is performed on the **Global Index** using the current task's dataset vector ( $D_{\text{vec}}$ ). This step efficiently identifies a small set of candidate datasets (e.g., the top 5) that are structurally similar to the current one, effectively pruning the search space.

<u>Fine Retrieval</u>: Next, a precise search is conducted on the **Step-wise Index**. The search is restricted only to the entries belonging to the candidate datasets identified in the coarse step. An exact k-nearest neighbor search (using cosine similarity) is performed with the current state vector  $s_{\text{vec}}$  to find the top k historical states that most closely match the current situation. The transitions ( $s_{\text{vec}}, a_t, r_t, s'_{\text{vec}}$ ) associated with these states are then returned to inform the LLM's prompt construction.

# A.7 Prompt Construction and Pipeline Generation

**Context-Aware Prompt Construction** At each invocation, the LLM Policy Advisor constructs a comprehensive prompt tailored to the current state  $s_t$ . This dynamic prompt is not redefined but is adapted based on the dataset's characteristics and the history of applied operations. The construction process follows these steps:

<u>State Encoding:</u> The current state  $s_t$ , which captures both the statistical and semantic features of the dataset  $D_t$ , is prepared for analysis. The statistical features include metrics such as mean, variance, and distribution characteristics, while the semantic features encode the task domain and the relationships between features.

Experience-Enhanced Context: The vectorized representation of the current state  $s_t$  is used to query the Experience Pool. This retrieval step aims to find the most relevant historical examples (i.e., previous dataset states that are similar to the current one). These historical examples consist of past actions, rewards, and subsequent states, which provide rich contextual information for the LLM.

**LLM-Guided Pipeline Generation** Using the constructed prompt, the LLM processes the context and generates a list of proposed pipeline. The LLM does not simply output individual actions (preprocessing operators) but instead suggests complete or partial pipelines. Each propsed pipeline  $\mathcal{P}_i$  is accompanied by:

- A confidence score  $c_i \in [0,1]$ , which indicates the LLM's certainty about the effectiveness of the suggested pipeline.
- A rationale r<sub>i</sub>, which is textual explanation provided by the LLM, describing why the proposed pipeline is well-suited for the current dataset.

The output is formalized as:

$$\mathcal{P}_{\text{suggested}} = \{ (P_i, c_i, r_i) \}_{i=1}^m \tag{5}$$

where  $P_i$  is a pipeline,  $c_i$  is the confidence score, and  $r_i$  is the textual explanation. The pipelines are ranked by their confidence scores, and the highest-ranked ones are passed on to the next stage for selection.