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Paper under double-blind review

ABSTRACT

Effectively fusing scarce high-accuracy data with massive but noisy low-accuracy data is a common challenge faced by machine learning across various fields, including agriculture, medicine, and remote sensing. Existing methods, which either directly concatenate datasets while ignoring accuracy differences or employ static weighting for training, struggle to achieve optimal performance. To address this, we introduce a deep learning framework incorporating a dynamic discard mechanism (DDL) that manages mixed-accuracy data through the selective, dynamic removal of low-accuracy instances characterized by high Mean Absolute Error (MAE) and the application of an adaptive weighting scheme. Our study validated this approach using rice cultivation data from China's four major rice-growing regions: South, Central, North, and Northeast China. Using site characteristics and nitrogen application rates as feature variables and rice yield as the target variable, we designated the high-accuracy dataset as the test set. Compared to machine learning models that process only single-accuracy datasets and other models designed for mixed-accuracy data, our DDL framework demonstrated a performance improvement of over 10% in metrics such as RMSE, MAE, and MAPE, achieving significantly higher prediction accuracy. A crop yield prediction model capable of handling multiple datasets simultaneously holds significant practical value for policymakers and other stakeholders. The dynamic discard mechanism and adaptive weighting algorithm employed by DDL also have considerable reference value for applications in other domains.

1 INTRODUCTION

Accurate estimation of crop yields is crucial for ensuring global food security and maintaining a stable market economy (Basso & Liu, 2019; Lecerf et al., 2019). This is fundamental for developing reasonable agricultural policies and effective food management (Zambrano et al., 2018), and also promotes sustainable agricultural development. In genuine agricultural contexts, crop productivity is synergistically governed by an intricate confluence of diverse parameters, notably edaphic and climatic factors, as well as nutrient application rates (Lai et al., 2024; Shuai & Basso, 2022). The intricate nature of these environmental variables makes it challenging to acquire a sufficient amount of high-accuracy data, which constrains the improvement of model prediction accuracy and generalization ability. In this context, building a prediction model that can effectively integrate datasets of mixed accuracy has become the core approach to enhancing estimation accuracy and generalization capability.

Agricultural data obtained directly from field observations and experiments are highly accurate and reliable but are limited in sample size due to high costs and labor-intensive processes. While existing research has used various machine learning (ML) models and deep neural networks (DNNs) to estimate crop yields and improve prediction accuracy (Akkem et al., 2023; Han et al., 2025). While most of these studies rely on data augmentation to process existing datasets for model training. This paucity of adequate sample diversity invariably compromises model generalizability, thereby failing to ameliorate the intrinsic data scarcity bottleneck inherent in agricultural domains.

In contrast, data generated through existing agricultural process models offer a massive sample size but are of lower accuracy. The simultaneous existence of high- and low-accuracy data presents both a significant challenge and a potential solution to agricultural data scarcity. However, the effective

054 fusion of scarce high-accuracy data with vast but noisy low-accuracy data remains an unresolved
 055 problem. Most existing studies either directly concatenate datasets, thereby overlooking disparities
 056 in data accuracy, and attempt to mitigate the noise introduced by lower-accuracy data through
 057 preprocessing methods (Zhang et al., 2022a). However, these approaches struggle to suppress the
 058 negative impact on model training while simultaneously preserving the intrinsic information content
 059 of the data. Alternatively, some studies employ static weighting strategies to enhance the fusion
 060 effect of mixed-accuracy data (Gao & Xie, 2025), yet these methods inherently lack adaptability to
 061 the dynamic changes occurring throughout the model training process.

062 To address these issues, this study proposes and validates a deep learning framework based on a
 063 dynamic discard mechanism (DDL), which can train datasets of different accuracies simultaneously.
 064 By leveraging a dynamic discard algorithm and an adaptive weighting mechanism, the framework
 065 enhances the model’s prediction accuracy and generalization ability, effectively solving the mixed-
 066 accuracy fusion problem.

067 The remainder of this paper is organized as follows: Section 2 reviews related work in this field;
 068 Section 3 describes the data acquisition strategies and sources; Section 4 presents the DDL archi-
 069 tectural design and various mechanisms; Section 5 shows the model results; and Section 6 discusses
 070 the research findings and outlines future research directions.

072 2 RELATED WORK

074 Recent research has increasingly focused on strategies to address the challenges of integrating het-
 075 erogeneous or uncertain datasets into predictive modeling.

077 **CNN–GAN–based methods** Methods combining Convolutional Neural Networks (CNNs) and
 078 Generative Adversarial Networks (GANs) have been widely used to improve data quality and aug-
 079 ment training samples. By learning high-level feature representations and generating realistic syn-
 080 thetic samples, these approaches effectively mitigate the issue of insufficient training data and en-
 081 hance model generalization. However, they typically assume a homogeneous data accuracy, often
 082 performing poorly when there are systematic differences in the reliability of training samples(Zhang
 083 et al., 2022b).

084 **U-Net with ConvLSTM architectures** To capture spatiotemporal dependencies in agricultural
 085 and environmental applications, researchers have proposed hybrid models that fuse the U-Net archi-
 086 tecture with Convolutional Long Short-Term Memory (ConvLSTM) modules. These models effec-
 087 tively integrate sequential information with spatial context, leading to significant improvements in
 088 tasks like crop monitoring. Nevertheless, their fusion process relies on a static architecture, lacking
 089 a dynamic mechanism to adapt to variations in input data quality(Kamangir et al., 2025).

090 **Remote sensing data assimilation with SCE-UA** Another research path involves data assimila-
 091 tion methods, such as utilizing optimization algorithms like the Shuffled Complex Evolution Al-
 092 gorithm (SCE-UA) to combine remote sensing observations with process-driven models. These
 093 approaches explicitly merge observations and simulated values, and can significantly enhance pre-
 094 dictive accuracy, particularly when observations are sparse or noisy. However, they typically op-
 095 erate within a deterministic optimization framework, lacking a mechanism to adaptively discard or
 096 re-weight low-accuracy samples during training(Li et al., 2024).

097 **Dynamic Reweighting and Data Selection** While our work shares the goal of improving training
 098 dynamics with methods like Population Based Augmentation (PBA) (Ho et al., 2019) and Sam-
 099 ple Reweighting (Ren et al., 2018), there is a fundamental distinction necessitated by the nature
 100 of mixed-accuracy scientific data. Traditional reweighting methods assign soft weights to high-loss
 101 samples, effectively down-weighting outliers but retaining them in the optimization process. In the
 102 context of simulations, high-error samples often represent systematic failures rather than aleatoric
 103 noise. Retaining these samples, even with low weights, risks corrupting the feature manifold.

104 In contrast, DDL employs a hard Dynamic Discard mechanism. By completely removing samples
 105 that persistently diverge from the high-accuracy distribution, DDL prevents the model from fitting
 106 to systematic biases, offering a more robust solution for integrating heterogeneous scientific datasets
 107 than augmentation or static reweighting strategies.

108 **3 DATA STRATEGY AND SOURCES**
 109

110 The primary objective of this study is to enhance the predictive accuracy and generalization capa-
 111 bility of the model to the greatest extent possible. However, during data collection, we identified
 112 a pervasive challenge: the quality and quantity of the required data exhibit an inverse relationship,
 113 resulting in two distinct types of data sources. To address this challenge, we propose an integrative
 114 strategy that combines the use of both high-accuracy and low-accuracy data.

115 The central premise of our data strategy is to combine the complementary strengths of these two
 116 sources—namely, the precision of high-accuracy data and the breadth of low-accuracy data—to
 117 construct a more robust and high-performing predictive model.

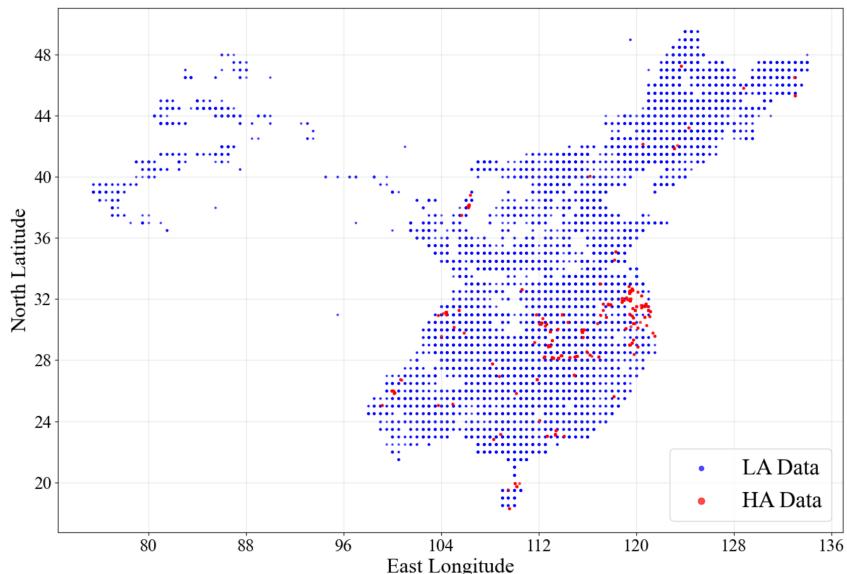
119 **3.1 HIGH-ACCURACY DATA: COLLECTED EMPIRICAL OBSERVATIONS**
 120

121 We established a high-accuracy, field-scale rice yield dataset through systematic field observations
 122 conducted between 2021 and 2024 across four major rice-growing regions. All data were collected
 123 under standardized experimental protocols with rigorous quality control procedures, ensuring con-
 124 sistency and reliability. In total, 704 observational records were obtained, providing a comprehen-
 125 sive empirical basis for subsequent modeling and analysis.

127 **3.2 LOW-ACCURACY DATA: DNDC-BASED SIMULATIONS**
 128

129 The DNDC (DeNitrification-DeComposition) model (Li et al., 1992) is a process-based biogeo-
 130 chemical model of carbon and nitrogen dynamics in agroecosystems. By coupling microbial
 131 metabolic processes with the soil’s physical environment, DNDC enables refined simulations of
 132 C–N cycles in complex agricultural systems. In this study, we employed version 9.5 of the DNDC
 133 model to simulate crop planting from paddy fields.

134 Soil property data and climate data obtained from the National Meteorological Science Data Center
 135 (2024) were aggregated at a 0.5° resolution into a format compatible with DNDC input requirements.
 136 We then ran the DNDC model in Region Mode, the simulation outputs include process-level crop
 137 growth data. From these results, we selected key variables including latitude/longitude, SOC, clay
 138 content, pH, bulk density (BD), average temperature, precipitation, irrigation, nitrogen application,
 139 crop type, and yield. In total, we obtained 43447 simulated records.



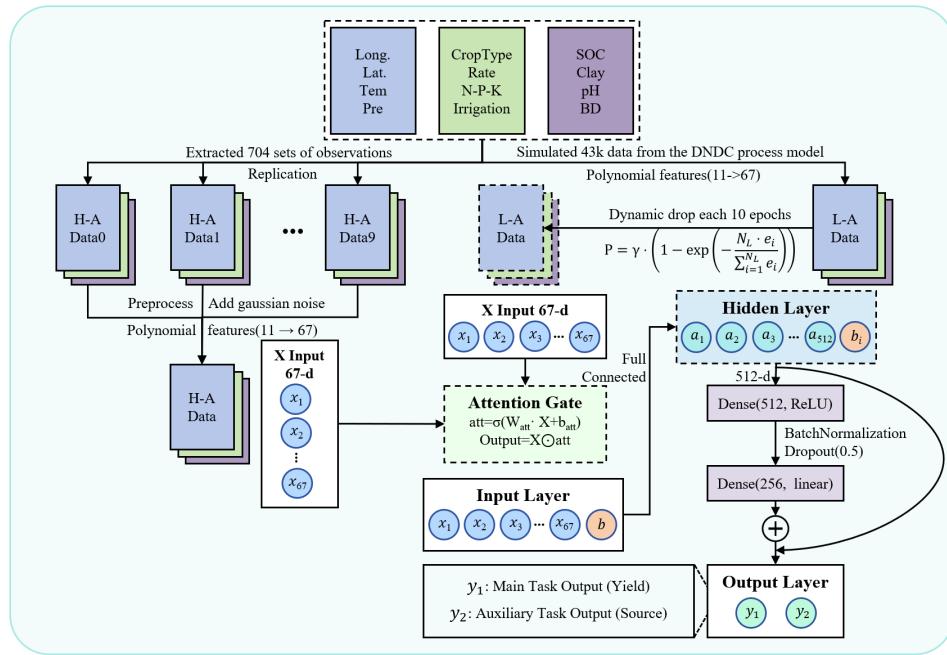
160 **Figure 1: Mixed-accuracy datasets distribution**
 161

162 4 METHODS
163164 4.1 NOTATION AND PROBLEM SETTING
165

166 This study addresses the problem of joint prediction by leveraging a small-sample, high-accuracy
167 dataset and a large-sample, low-accuracy dataset. Let $D^H = \{(x_i^H, y_i^H)\}_{i=1}^{n_H}$ and $D^L =$
168 $\{(x_i^L, y_i^L)\}_{i=1}^{n_L}$ denote the high-accuracy (collected) data and the low-accuracy (simulated) data re-
169 spectively. Here, x_i represents an 11-dimensional feature vector: **Longitude**, **Latitude**, **Tem** (tem-
170 **perature**), **Pre** (precipitation), **rate** (fertilizer application **N** rate), **SOC** (soil organic carbon), **Clay**
171 (**clay content**), **pH** (soil pH), **BD** (soil bulk density), **irrigation**, and **CropType** and y_i is the target
172 variable, **Yield** (rice yield per mu). The dataset sizes are $n_H = 704$ and $n_L = 43,447$. Our primary
173 objective is to achieve superior predictive accuracy on a test set partitioned from the high-accuracy
174 dataset, while also maximizing the model’s generalization capability. **In our pipeline the high-ac-**
175 **curacy dataset was split prior to model training: 20% of D^H were reserved as a final, held-out test**
176 **set, 20% of the remaining high-accuracy samples were used as a validation set, and the remainder**
177 **were used for training (all random splits used seed 42).** The two datasets therefore share the same
178 **feature / label schema but differ in measurement accuracy and noise characteristics.** The held-out
179 **D^H test subset was not used in any way during training, dynamic discard procedures, or adaptive**
180 **weight tuning.**

182 4.2 ARCHITECTURE AND TECHNICAL DESIGN OF THE DDL FRAMEWORK
183

184 As illustrated in Figure 1, we propose the DDL architecture, which consists of an input layer, a
185 **Feature Attention Gating Module**, a **Modified Residual Block**, a fully connected layer, and a
186 multi-task output layer. The model incorporates a **dynamic dropout mechanism** and **adaptive**
187 **dynamic weights**. The specific technical principles of these components are detailed below.



211 Figure 2: **Schematic overview of the DDL framework.** The architecture comprises a Feature Atten-
212 **Gation Module for input reweighting, Modified Residual Blocks to mitigate gradient vanishing,**
213 **and a multi-task output layer. The Dynamic Discard mechanism and Adaptive Dynamic Weights**
214 **operate during training to filter noise and balance task losses.**

216 4.2.1 DYNAMIC DISCARD STRATEGY
217

218 The core mechanism of our model is the **Dynamic Discard (DD) strategy**, which is not a static
219 data cleaning process performed before training. Instead, it is a progressive filtering mechanism that
220 evaluates and discards low-accuracy data samples during the training process based on the model’s
221 current predictive performance, thereby gradually filtering out noisy data.

222 Specifically, at epoch t , with current model parameters θ_t , we first perform a forward pass on all
223 low-accuracy data samples in $D^L = \{(x_i^L, y_i^L)\}_{i=1}^{N_L}$ to obtain their primary task (yield prediction)
224 predictions, $\hat{y}_i^L = f_{\theta_t}(x_i^L)$. We then calculate the absolute prediction error $e_i = |\hat{y}_i^L - y_i^L|$ for each
225 low-accuracy sample.

226 To ensure the discard threshold adapts to the model’s performance at different training stages, we
227 compute the Mean Absolute Error (MAE) (Hodson, 2022) of all low-accuracy samples in the current
228 batch as a dynamic baseline:

$$229 MAE_t = \frac{1}{n_L} \sum_{i=1}^{n_L} |\hat{y}_i^L - y_i^L|$$

230 Next, we normalize each sample’s error to obtain a deviation metric relative to the model’s current
231 average performance: $\tilde{e}_i = e_i / MAE_t$.

232 Based on this deviation, we calculate a dynamic discard probability P_i for each sample, determined
233 by an exponential decay function:

$$234 P_i = \gamma \cdot (1 - \exp(-\tilde{e}_i))$$

235 where γ is a base discard probability factor that controls the steepness of the probability curve,
236 thereby regulating the penalty on high-error samples. Finally, a Bernoulli sampling process (Yu
237 et al., 2022) determines whether the sample is kept: a random number r is generated in the range
238 $[0, 1]$. If $r < P_i$, the sample is discarded; otherwise, it is retained for the current and subsequent
239 gradient updates.

240 This process can be conceptualized as a filtering operator \mathcal{D}_t that acts on the low-accuracy dataset,
241 outputting a filtered subset:

$$242 \mathcal{D}_t(\{(x_i^L, y_i^L)\}) = \{(x_j^L, y_j^L) \mid r_j \geq p_j\}$$

243 This subset is then combined with the high-accuracy data D^H to form the training data for epoch
244 t . Through this mechanism, the model can initially leverage the large volume of low-accuracy data
245 to quickly learn general features. With advancing training and enhanced model performance, the
246 DD mechanism adopts a more stringent approach, systematically filtering out low-quality samples
247 that persistently yield high errors and are likely indicative of noise or substantial divergence from
248 the true distribution. This allows the model to later focus on refining its predictive capabilities using
249 higher-quality data, effectively preventing the negative influence of low-quality data and achieving
250 a dynamic balance between data quality and quantity.

251 In summary, the dynamic discard probability formula of the proposed model is:

$$252 P_i = \gamma \cdot \left(1 - \exp \left(\frac{-N_L \cdot e_i}{\sum_{j=1}^{N_L} e_j} \right) \right)$$

253 4.2.2 FEATURE ATTENTION GATING MODULE AND REGULARIZATION
254

255 The model’s input layer receives 11 types of features, including soil parameters, meteorological
256 factors, and crop data. To enhance the importance of features highly correlated with the target variable,
257 we designed a trainable **Feature Attention Gating Module** (Meng et al., 2022; Dhingra et al.,
258 2016). This module uses fully connected layers (Basha et al., 2020) to perform dynamic feature
259 weighting at the input layer, unlike the more computationally expensive self-attention mechanism,
260 ensuring core features play a primary role in the final prediction.

261 The mathematical formulation is as follows:

$$262 \text{att} = \sigma(W_{\text{att}} \cdot x + b_{\text{att}})$$

output = $x \odot \text{att}$

where σ is the attention gate vector, W_{att} and b_{att} are trainable parameters, and \odot denotes element-wise multiplication.

To prevent model overfitting, we introduce L2 regularization (Van Laarhoven, 2017), which modifies the total objective function to:

$$\mathcal{L}_{req} = \mathcal{L} + \lambda \left\| W_{att} \right\|_F^2$$

where $\lambda = 10^{-4}$ is the regularization coefficient and $\|\cdot\|_F$ is the Frobenius norm of the weight matrices. This regularization term can be interpreted as a Gaussian prior on the weight parameters, ensuring that the attention weights do not become overly concentrated on a few features.

4.2.3 MODIFIED RESIDUAL BLOCK

To address the issue of **gradient vanishing** in deep networks (Tan & Lim, 2019), we designed a modified residual block structure (Zhang et al., 2017; Tang et al., 2024). Each block consists of two fully connected layers: the first employs a **ReLU** activation function (He et al., 2018) to introduce non-linearity, while the second uses a linear activation to maintain numerical stability. A skip connection then adds the block's input to its output.

The mathematical expression is as follows:

$$x^{(l+1)} = g\left(x^{(l)}\right) + \mathcal{F}\left(x^{(l)}\right)$$

where $g(x^{(l)})$ is a dimension-adapting function that acts as an identity mapping when the input dimension $\dim(x^{(l)})$ matches the target output dimension d_{out} ; otherwise, it performs a dimensionality transformation via a projection matrix W_s :

$$g\left(x^{(l)}\right) = \begin{cases} x^{(l)}, & \text{if } \dim(x^{(l)}) = d_{\text{out}} \\ W_s x^{(l)} + b_s, & \text{otherwise} \end{cases}$$

The residual function $\mathcal{F}(x^{(l)})$ is implemented with two fully connected layers, and its computational flow is given by the following equation:

$$\mathcal{F}(x^{(l)}) = W_2^{(l)} \cdot z_3 + b_2^{(l)}$$

where z_3 is an intermediate variable computed as follows:

$$z_1 = \delta_{\text{ReLU}} \left(W_1^{(l)} x^{(l)} + b_1^{(l)} \right)$$

$$z_2 = \text{BN}(z_1)$$

$$z_3 = \text{Dropout}(z_2)$$

Here, δ_{ReLU} denotes the ReLU activation function. We constructed these identity mapping paths within the 512- and 256-dimensional hidden layers. This allows gradients to be passed directly to the shallower layers during backpropagation, effectively mitigating the vanishing gradient problem. Combined with **Dropout**, this structure also provides a strong regularization effect.

4.2.4 ADAPTIVE DYNAMIC WEIGHTS

To enable the model to learn high-accuracy predictions from the small dataset while simultaneously improving its generalization from the large dataset, we designed a framework with **adaptive dynamic weights** (Yang et al., 2022; Xiao & Zhang, 2021; Cao et al., 2023). This approach dynamically balances the loss weights between the main task (regression prediction) and an auxiliary task (data source classification).

The multi-task learning framework consists of a main task (predicting the target variable via regression) and an auxiliary task (predicting the data source as a binary classification task). The overall loss function is defined as:

$$f_{total} \equiv \alpha \cdot f_{main} + (1 - \alpha) \cdot f_{aux}$$

324 where α is a dynamic weighting coefficient, $\mathcal{L}_{\text{main}}$ is the mean squared error (MSE) loss, and \mathcal{L}_{aux} is
 325 the classification cross-entropy loss:

$$327 \quad L_{\text{main}} = \frac{1}{N} \sum_{i=1}^N (y_i - \hat{y}_i)^2$$

$$330 \quad \mathcal{L}_{\text{aux}} = -\frac{1}{N} \sum_{i=1}^N [s_i \log(\hat{s}_i) + (1 - s_i) \log(1 - \hat{s}_i)]$$

333 The initial value of α is set to 0.4 and is increased exponentially with each epoch, up to a specified
 334 limit. This mechanism ensures that the model initially leverages the classification task to enhance
 335 its generalization, then later focuses on optimizing predictive accuracy as the main task's influence
 336 increases. The dynamic weight update mechanism is as follows:

$$337 \quad \alpha_{\text{epoch}+1} = \max(0.4, \min(0.8, \alpha_{\text{epoch}} \times 1.005))$$

339 5 RESULT

341 5.1 EXPERIMENTAL RESULTS AND DISCUSSION

343 The experimental results demonstrate that the proposed DDL framework significantly outperforms
 344 existing mainstream methods for the rice yield prediction task. As shown in Table 1, DDL exhibits
 345 superior performance across all four metrics: MAE, RMSE, R², and MAPE, surpassing traditional
 346 machine learning models (e.g., XGBoost, Random Forest), process-based models (e.g., DNDC), and
 347 domain-specific hybrid architectures (e.g., CNN+GAN, Remote Sensing Assimilation + SCE-UA).
 348 The framework achieves an approximate 10% improvement on these metrics, with the predictive R²
 349 value reaching 0.68. This strong agreement between predicted and observed yields is further visual-
 350 ized in the scatter plot of Figure 3, where most samples cluster closely around the fitted regression
 351 line, particularly in the mid-to-high yield range.

352 Table 1: Performance Comparison of Various Models on Regression Tasks

354 Model	MAE ↓	RMSE ↓	R ² ↑	MAPE ↓	Reference
355 ML Model					
356 Gradient Boosting R	1001.96	1287.90	0.6348	0.1648	(Friedman, 2001)
357 LightGBM	1011.40	1316.29	0.6186	0.1678	(Ke et al., 2017)
358 Random Forest	1076.34	1436.98	0.5454	0.1754	(Breiman, 2001)
359 XGBoost Regression	1064.40	1388.34	0.5757	0.1750	(Chen & Guestrin, 2016)
360 GPR	1296.37	1808.19	0.3381	0.2098	(Rasmussen & Williams, 2006)
MHA-MLP	1115.07	1496.70	0.5465	0.1767	-
361 Related Work					
362 CNN+GAN	1115.56	1637.84	0.3587	0.2026	(Zhang et al., 2022b)
363 UNet-ConvLSTM	1075.91	1357.26	0.5497	0.1672	(Kamangir et al., 2025)
364 PBA-ResNet	1082.35	1462.31	0.4823	0.1612	(He et al., 2016)
365 PBA-MLP	1108.91	1642.98	0.4535	0.1635	(Ho et al., 2019)
Broad Learning System	1160.32	1559.85	0.5074	0.1815	(Liu & Chen, 2018)
366 Process Model					
367 DNDC Model	1439.35	1754.55	0.2910	8.7361	(Li et al., 1992)
368 DDL Model					
369 Ours	852.30	1212.14	0.6837	0.1424	-

371 5.2 ABLATION STUDY

373 An ablation study was conducted to validate the necessity and effectiveness of each component
 374 within the DDL framework. The results confirm that the absence of any single component leads
 375 to a decline in model performance. Specifically, a "single-accuracy" version of the model trained
 376 exclusively on the high-accuracy data demonstrated significantly limited performance. Similarly,
 377 models that did not employ the **dynamic discard mechanism** during mixed-precision training or
 those that used a static weighting strategy in place of the **adaptive dynamic weights** both performed

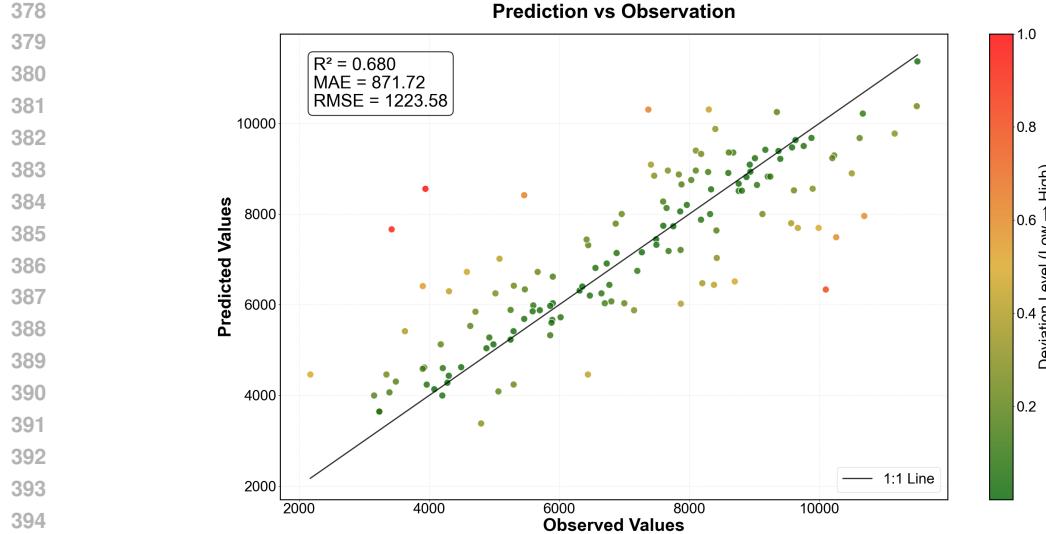


Figure 3: Scatter plot of predicted versus observed rice yield values. Each point corresponds to a sample, with its color indicating the absolute deviation from the fitted regression line: green denotes close agreement (small error), gradually transitioning to red for larger deviations (poor agreement)

worse than the complete framework. These findings in Table 2 suggest that static data fusion or the unfiltered use of low-accuracy data fails to effectively suppress noise interference. The synergistic design of the **dynamic discard** and **adaptive dynamic weights** is, therefore, crucial for DDL’s ability to achieve high-precision predictions. The results of the ablation study fully validate the necessity and efficacy of the proposed mechanisms in fusing mixed-precision data.

Table 2: Ablation Study of the DDL Framework

Model	MAE ↓	RMSE ↓	R ² ↑	MAPE ↓
ML Model				
Gradient Boosting R	1001.96	1287.90	0.6348	0.1648
Related Work				
UNet-ConvLSTM	1075.91	1357.26	0.5497	0.1672
Process Model				
DNDC Model	1439.35	1754.55	0.2910	8.7361
DDL Model				
Ours(High-Accuracy Only)	958.62	1335.92	0.6158	0.1602
Ours(w/o Dynamic Discard)	887.10	1273.89	0.6506	0.1477
Ours(w/ Static Weighting)	870.31	1249.90	0.6637	0.1439
Ours(Full)	852.30	1212.14	0.6837	0.1424

5.3 GENERALIZATION ABILITY

Cross-Species Transfer To assess the generalization ability of our proposed DDL approach beyond rice, we conducted cross-species experiments on Wheat and Maize. For Wheat we further collected a high-accuracy dataset of 500 samples and a low-accuracy dataset of 52,450 samples; for Maize the high-accuracy dataset contains 567 samples and the low-accuracy dataset contains 38,492 samples. We retrained and evaluated our full DDL model on each species and compared its predictive performance against a broad set of baselines. Tables 8 and 9 summarize the results for each method. From the results shown in these tables, it’s evident that—even when evaluated on other crops—our DDL model remains highly competitive.

Sparse-Region Generalization To evaluate model robustness across regions with varying data densities, we designed a validation experiment based on spatial grid partitioning. Specifically, the geographic space was divided into grids with a resolution of $1^\circ \times 1^\circ$ in both latitude and longitude,

432 Table 3: All Performance Comparison of Various Models on Regression Tasks for Wheat
433

434 Model	435 MAE ↓	436 RMSE ↓	437 R ² ↑	438 MAPE ↓	439 Reference
ML Models					
440 Gradient Boosting R	441 1143.22	442 1576.82	443 0.4907	444 0.2698	445 (Friedman, 2001)
446 LightGBM	447 1061.73	448 1445.15	449 0.5722	450 0.2651	451 (Ke et al., 2017)
452 MHA-MLP	453 1200.85	454 1579.87	455 0.4802	456 0.2707	457 -
Related Work					
458 PBA-MLP	459 1134.83	460 1508.56	461 0.4750	462 0.2638	463 (Ho et al., 2019)
464 Broad Learning System	465 1197.73	466 1613.30	467 0.4580	468 0.2839	469 (Liu & Chen, 2018)
DDL Model					
470 Ours(Full)	471 1118.22	472 1434.27	473 0.5590	474 0.2641	475 -

445 Table 4: All Performance Comparison of Various Models on Regression Tasks for Maize
446

447 Model	448 MAE ↓	449 RMSE ↓	450 R ² ↑	451 MAPE ↓	452 Reference
ML Models					
453 Gradient Boosting R	454 1513.57	455 2013.99	456 0.6838	457 0.3206	458 (Friedman, 2001)
459 LightGBM	460 1506.19	461 2032.21	462 0.6781	463 0.3251	464 (Ke et al., 2017)
465 Random Forest	466 1561.46	467 2063.54	468 0.6681	469 0.3350	470 (Breiman, 2001)
Related Work					
471 PBA-MLP	472 2115.14	473 2706.12	474 0.3155	475 0.6853	476 (Ho et al., 2019)
477 Broad Learning System	478 1672.19	479 2335.77	480 0.3393	481 0.2178	482 (Liu & Chen, 2018)
DDL Model					
483 Ours (Full)	484 1382.21	485 1847.46	486 0.6735	487 0.2542	488 -

459 and the density of both high- and low-accuracy data within each grid was quantified. Based on these
460 densities, we constructed different testing scenarios (Table ??).
461

462 As shown in the results, reducing the quantity of high-accuracy data (transitioning from Case A
463 to Case B) results in a moderate performance decrease, with R^2 dropping from 0.7480 to 0.7115.
464 While this indicates that high-accuracy data contributes to refining predictions, the model main-
465 tains a relatively high performance level. In stark contrast, limiting the availability of low-accuracy
466 data (Case C) leads to a substantial deterioration in performance (R^2 plummets to 0.2644), even
467 when high-accuracy data is abundant. These findings refute the notion that high-accuracy data alone
468 is sufficient for broad generalization and underscore that the massive low-accuracy dataset acts as
469 a critical stabilizer, enabling the DDL framework to generalize effectively even in regions where
470 empirical observations are sparse.
471

472 Table 5: Generalization Ability of the DDL Framework
473

474 Data Configuration	475 MAE ↓	476 RMSE ↓	477 R ² ↑	478 MAPE ↓
479 Case A	480 810.27	481 1091.62	482 0.7480	483 0.1308
484 Case B	485 616.35	486 867.91	487 0.7115	488 0.0835
489 Case C	490 1386.74	491 1944.45	492 0.2644	493 0.2649
494 Case D	495 -	496 -	497 -	498 -

499 **Note:** Grids were ranked by the **density of high-accuracy samples**.

500 - **Top 50%**: Regions with high density of empirical observations (Data-Rich).

501 - **Bottom 50%**: Regions with low density of empirical observations (Data-Sparse).

502 *Configuration Details:*

503 - **Case A**: Full low-accuracy data + High-accuracy data from Top 50% regions.

504 - **Case B**: Full low-accuracy data + High-accuracy data from Bottom 50% regions.

505 - **Case C**: Sparse low-accuracy data + High-accuracy data from Top 50% regions.

506

486 5.4 SENSITIVITY TO NOISE IN LOW-ACCURACY DATA
487488 To test whether the predictive accuracy of our DDL model is tightly constrained by the quality of the
489 low-accuracy dataset, we injected zero-mean Gaussian noise into the low-accuracy data at relative
490 magnitudes of 0%, 10%, 20%, 30% and 40% (noise standard deviation expressed as a fraction of the
491 original signal standard deviation). We retrained and evaluated the full DDL model under each noise
492 condition and compared its R^2 against representative baselines. Results in Table 6 show that even
493 with 20% Gaussian noise the DDL model outperforms the strongest baselines, indicating limited
494 sensitivity to simulation noise in the low-accuracy source.
495496 Table 6: Robustness to Noise Injected into Low-Accuracy Data (R^2)
497

498 Model / Condition	499 $R^2 \uparrow$	500 Reference
501 Machine-Learning Baselines		
502 Gradient Boosting Regression	0.6348	(Friedman, 2001)
503 LightGBM	0.6186	(Ke et al., 2017)
504 XGBoost Regression	0.5757	(Chen & Guestrin, 2016)
505 Related Work		
506 UNet-ConvLSTM	0.5497	(Kamangir et al., 2025)
507 Broad Learning System	0.5074	(Liu & Chen, 2018)
508 DDL (Ours) — noise levels applied to low-accuracy data		
509 Ours (0% noise)	510 0.6837	-
511 Ours (10% noise)	0.6681	-
512 Ours (20% noise)	0.6452	-
513 Ours (30% noise)	0.6134	-
514 Ours (40% noise)	0.5662	-

515 6 CONCLUSION AND FUTURE WORK
516517 This study proposes a Deep Learning framework with Dynamic Dropping mechanism, designed to
518 achieve a deep and organic integration of large-scale low-accuracy datasets with small-scale high-
519 accuracy datasets. The framework provides an effective solution for further improving both the
520 accuracy and generalization ability of crop yield prediction models. At its core, DDL introduces a
521 dynamic dropping strategy, which contrasts with traditional static data-cleaning approaches that dis-
522 card noisy samples or retain uninformative data prior to training (Ashfaq et al., 2025). By embedding
523 data handling as a continuous, dynamic process throughout training, DDL enhances generalization
524 and stability on high-accuracy test sets. This strategy mitigates the inherent trade-off between data
525 quantity and data quality, reducing the reliance on stringent requirements for high-quality datasets.
526 Consequently, training with heterogeneous data sources becomes more reliable in domains such as
527 agriculture and industry, where access to high-quality data is limited (Paudel et al., 2022).528 Despite these promising results, three directions remain for further exploration. Firstly, the dropping
529 strategy in this work primarily relies on absolute prediction error, without accounting for intrinsic
530 data distribution characteristics. Future research may incorporate distributional differences between
531 datasets when estimating dropping probabilities (Egele et al., 2024). Secondly, the current weight
532 adjustment scheme follows a linear schedule. Although it alleviates the need for manual weight tun-
533 ing and provides some adaptivity, it lacks adjustments based on real-time training dynamics. Closing
534 this loop through validation-based feedback mechanisms would be critical for further improving pre-
535 dictive accuracy (Caljon et al., 2025). Eventually, while this study focuses on rice yield prediction
536 in agriculture, the proposed framework could be extended to other domains, offering a generalizable
537 solution for integrating heterogeneous data and enhancing model generalization (Zhang et al., 2025).538 In summary, the DDL framework presents a practical approach to jointly train on small high-ac-
539 curacy and large low-accuracy datasets; empirical results on crop yield tasks show consistent im-
540 provements over examined baselines. While these results are encouraging, we avoid broad claims
541 about universal generalizability and instead emphasize that DDL is a promising mechanism whose
542 applicability to other domains should be explored in future work.

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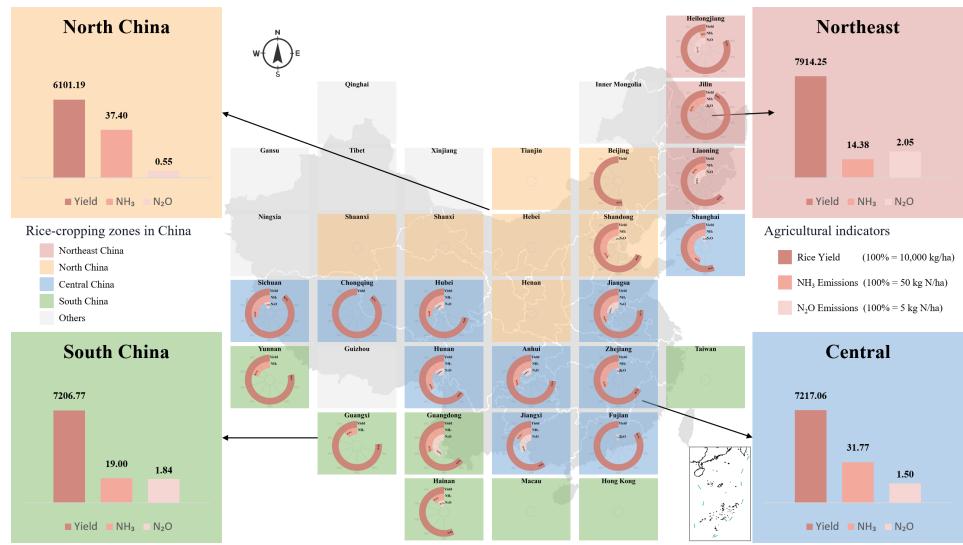
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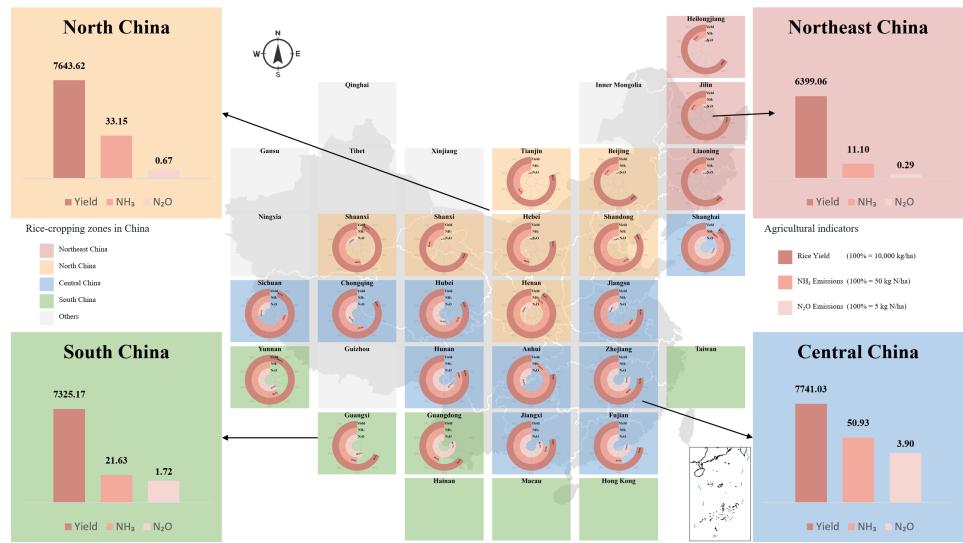
A APPENDIX

A.1 DATA DISTRIBUTION

The following shows the average yield per mu of the Mixed-accuracy dataset across four rice-growing regions and its distribution across various provinces, as illustrated in Figure 4a.

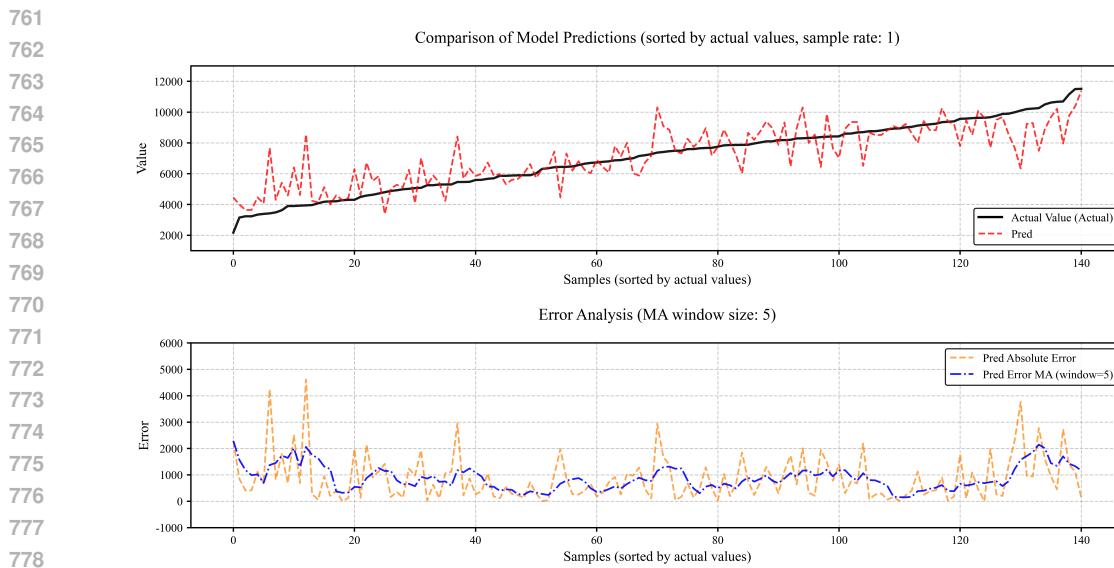
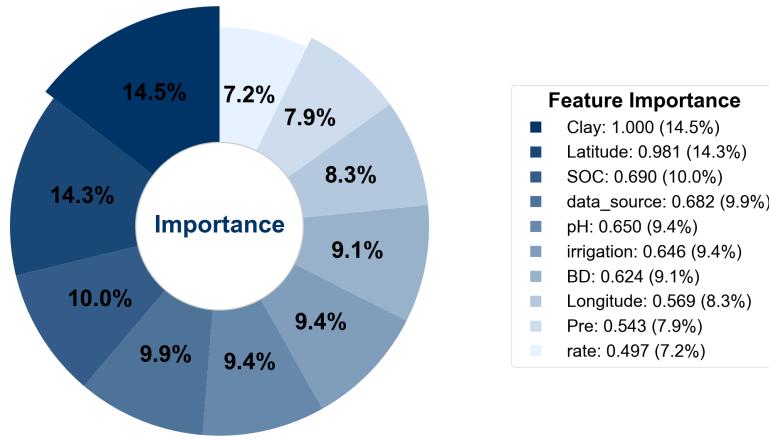


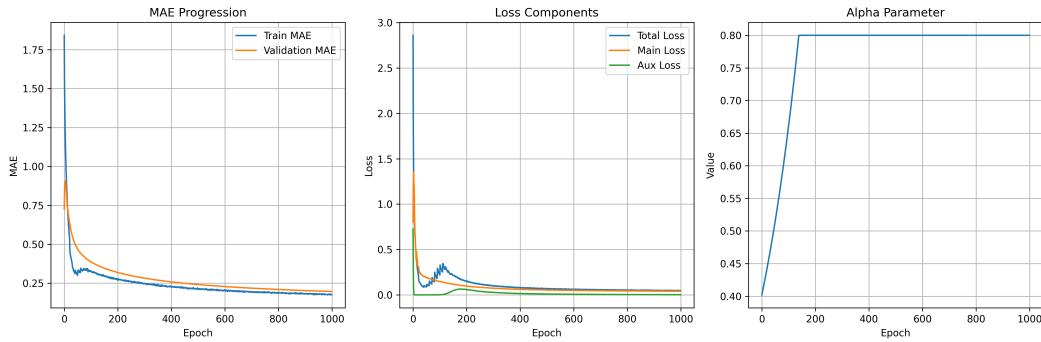
(a) Prediction performance on the Mixed-accuracy dataset (low-accuracy subset).



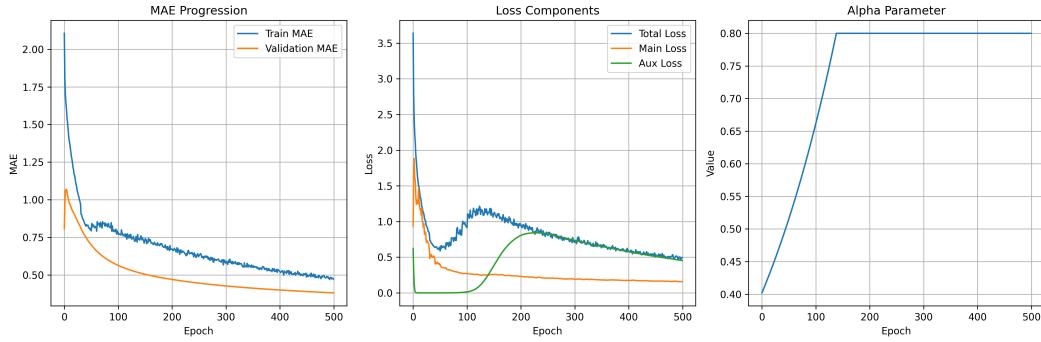
(b) Prediction performance on the Mixed-accuracy dataset (high-accuracy subset).

Figure 4: Comparison of model prediction accuracy across subsets of the Mixed-accuracy dataset.

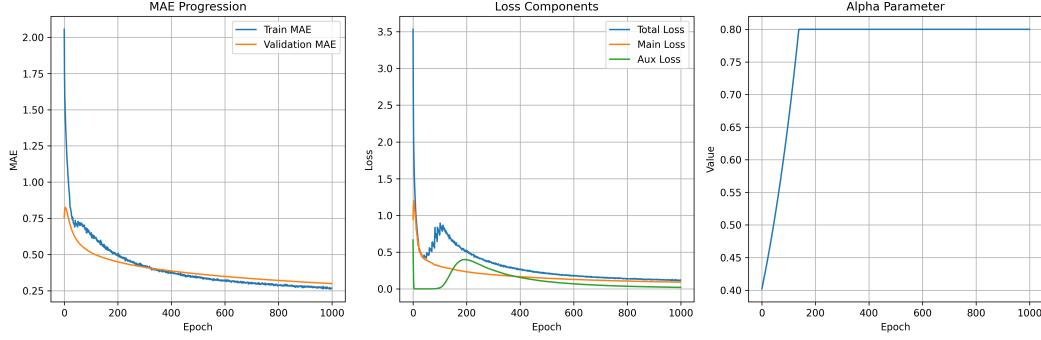
756 A.2 TRAINING DETAILS
757758 To further investigate model behavior, we present two complementary visualizations: one character-
759 izing prediction fidelity across samples, and another revealing the contribution of input features to
760 the model’s decisions.
761779 Figure 5: Temporal-style prediction and error analysis across samples. The upper panel displays the
780 actual rice yield (solid black line) alongside model predictions (dashed blue line) ordered by sample
781 index. The lower panel shows the per-sample absolute error (orange dashed line) and its moving
782 average (blue dash-dot line), computed using a sliding window to suppress noise and highlight
783 systematic bias or regional error patterns. A larger window width improves trend visibility but
784 reduces sensitivity to local error spikes.
785786 **Feature Importance Distribution**
787804 Figure 6: Relative importance of input features in the predictive model, visualized as a donut chart.
805 Features are ranked by their contribution to prediction accuracy, with the top- and bottom-ranked
806 features explicitly labeled and color-highlighted. The hollow center enhances visual focus on the
807 proportional influence of each variable, underscoring which agronomic or environmental factors
808 drive model performance.
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(a) Rice training curves



(b) Wheat training curves



(c) Maize training curves

Figure 7: Training curves of the DDL framework across three major crops. Each subplot shows three key metrics: (1) MAE progression for both training and validation sets, (2) Loss components including total loss, main loss (yield prediction), and auxiliary loss (data source classification), and (3) Alpha parameter evolution that dynamically balances the main and auxiliary tasks during training. The consistent convergence patterns across all three crops validate the generalizability of the DDL framework for different crop yield prediction scenarios.

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864 A.3 RESULT DETAILS
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866 Tables 7, 8, and 9 comprehensively evaluate the predictive performance of our DDL framework
867 against various baseline models across three major crops: rice, wheat, and maize. These tables
868 present detailed comparisons using multiple evaluation metrics. The results validate the general-
869 **izability of the DDL framework beyond rice to other important cereal crops.**

870
871 Table 7: All Performance Comparison of Various Models on Regression Tasks for Rice
872

873 Model	874 MAE ↓	875 RMSE ↓	876 R ² ↑	877 MAPE ↓	878 Reference
ML Model					
875 Decision Tree R	876 1098.62	877 1495.59	878 0.5076	879 0.1752	880 (Breiman et al., 1984)
876 Gradient Boosting R	877 1001.96	878 1287.90	879 0.6348	880 0.1648	881 (Friedman, 2001)
877 LightGBM	878 1011.40	879 1316.29	880 0.6186	881 0.1678	882 (Ke et al., 2017)
878 Linear Regression	879 1344.19	880 1684.27	881 0.3755	882 0.2309	883 (Seber & Lee, 2012)
879 Random Forest	880 1076.34	881 1436.98	882 0.5454	883 0.1754	884 (Breiman, 2001)
880 SVR	881 1777.09	882 2123.63	883 0.0072	884 0.3108	885 (Drucker et al., 1996)
881 XGBoost Regression	882 1064.40	883 1388.34	884 0.5757	885 0.1750	886 (Chen & Guestrin, 2016)
882 Deep Learning Regression	883 2938.72	884 3500.50	885 -1.6976	886 0.4183	887 (Goodfellow et al., 2016)
883 GPR	884 1296.37	885 1808.19	886 0.3381	887 0.2098	888 (Rasmussen & Williams, 2006)
884 MHA-MLP	885 1115.07	886 1496.70	887 0.5465	888 0.1767	889 -
Related Work					
885 CNN+GAN	886 1115.56	887 1637.84	888 0.3587	889 0.2026	890 (Zhang et al., 2022b)
886 UNet-ConvLSTM	887 1075.91	888 1357.26	889 0.5497	890 0.1672	891 (Kamangir et al., 2025)
887 Remote Sensing + SCE-UA	888 4025.99	889 5391.50	890 -2.7046	891 0.7144	892 (Li et al., 2024)
888 PBA-ResNet	889 1082.35	890 1462.31	891 0.4823	892 0.1612	893 (He et al., 2016)
889 PBA-MLP	890 1108.91	891 1642.98	892 0.4535	893 0.1635	894 (Ho et al., 2019)
890 Broad Learning System	891 1160.32	892 1559.85	893 0.5074	894 0.1815	895 (Liu & Chen, 2018)
Process Model					
891 DNDC Model	892 1439.35	893 1754.55	894 0.2910	895 8.7361	896 (Li et al., 1992)
DDL Model					
894 Ours(High-Accuracy Only)	895 958.62	896 1335.92	897 0.6158	898 0.1602	899 -
895 Ours(w/o Dynamic Discard)	896 887.10	897 1273.89	898 0.6506	899 0.1477	900 -
896 Ours(w/ Static Weighting)	897 870.31	898 1249.90	899 0.6637	900 0.1439	901 -
897 Ours(Full)	898 852.30	899 1212.14	900 0.6837	901 0.1424	902 -

900 Table 8: All Performance Comparison of Various Models on Regression Tasks for Wheat
901

902 Model	903 MAE ↓	904 RMSE ↓	905 R ² ↑	906 MAPE ↓	907 Reference
ML Model					
904 Decision Tree R	905 1326.32	906 1929.20	907 0.2377	908 0.3192	909 (Breiman et al., 1984)
905 Gradient Boosting R	906 1143.22	907 1576.82	908 0.4907	909 0.2698	910 (Friedman, 2001)
906 LightGBM	907 1061.73	908 1445.15	909 0.5722	910 0.2651	911 (Ke et al., 2017)
907 Linear Regression	908 1454.95	909 1817.08	910 0.3237	911 0.3653	912 (Seber & Lee, 2012)
908 Random Forest	909 1143.93	910 1617.93	911 0.4638	912 0.2849	913 (Breiman, 2001)
909 SVR	910 1819.56	911 2213.53	912 -0.0036	913 0.5500	914 (Drucker et al., 1996)
910 XGBoost Regression	911 1129.97	912 1634.62	913 0.4527	914 0.2770	915 (Chen & Guestrin, 2016)
911 Deep Learning Regression	912 2967.98	913 3651.12	914 -1.7305	915 0.5131	916 (Goodfellow et al., 2016)
912 GPR	913 1557.16	914 1945.91	915 0.2115	916 0.4009	917 (Rasmussen & Williams, 2006)
913 MHA-MLP	914 1200.85	915 1579.87	916 0.4802	917 0.2707	918 -
Related Work					
914 PBA-MLP	915 1134.83	916 1508.56	917 0.4750	918 0.2638	919 (Ho et al., 2019)
915 Broad Learning System	916 1197.73	917 1613.30	918 0.4580	919 0.2839	920 (Liu & Chen, 2018)
DDL Model					
917 Ours(Full)	918 1118.22	919 1434.27	920 0.5590	921 0.2641	922 -

918 Table 9: All Performance Comparison of Various Models on Regression Tasks for Maize
919

920 Model	921 MAE ↓	922 RMSE ↓	923 R ² ↑	924 MAPE ↓	925 Reference
ML Model					
926 Decision Tree R	927 1936.85	928 2685.83	929 0.4377	930 0.3633	(Breiman et al., 1984)
926 Gradient Boosting R	927 1513.57	928 2013.99	929 0.6838	930 0.3206	(Friedman, 2001)
926 LightGBM	927 1506.19	928 2032.21	929 0.6781	930 0.3251	(Ke et al., 2017)
926 Linear Regression	927 2215.98	928 2901.25	929 0.3439	930 0.4816	(Seber & Lee, 2012)
926 Random Forest	927 1561.46	928 2063.54	929 0.6681	930 0.3350	(Breiman, 2001)
926 SVR	927 2656.28	928 3589.16	929 -0.0042	930 0.6257	(Drucker et al., 1996)
926 XGBoost Regression	927 1635.27	928 2188.98	929 0.6265	930 0.3221	(Chen & Guestrin, 2016)
926 Deep Learning Regression	927 6155.25	928 6969.60	929 -2.7866	930 0.7119	(Goodfellow et al., 2016)
926 GPR	927 1901.65	928 2587.13	929 0.1895	930 0.2673	(Rasmussen & Williams, 2006)
926 MHA-MLP	927 1841.45	928 2493.63	929 0.2470	930 0.2619	-
Related Work					
932 PBA-MLP	933 2115.14	934 2706.12	935 0.3155	936 0.6853	(Ho et al., 2019)
932 Broad Learning System	933 1672.19	934 2335.77	935 0.3393	936 0.2178	(Liu & Chen, 2018)
DDL Model					
935 Ours(Full)	936 1382.21	937 1847.46	938 0.6735	939 0.2542	940 -

937 **B. ETHICS STATEMENT**

938 This research adheres to the ICLR Code of Ethics. No human subjects or animal experiments were
939 involved in this study. All datasets used were obtained in accordance with their respective usage
940 guidelines to ensure no violation of privacy. We have made every effort to avoid bias or discrimina-
941 tory outcomes in our research. No personally identifiable information was used, and no experiments
942 were conducted that could raise privacy or security concerns. We are committed to maintaining
943 transparency and integrity throughout this research.

944 **C. REPRODUCIBILITY STATEMENT**

945 We have made every effort to ensure that the results presented in this paper are reproducible. All code
946 and datasets have been submitted as supplementary materials to facilitate replication and verification
947 by others.

948 **D. LARGE LANGUAGE MODEL (LLM) USAGE STATEMENT**

949 A large language model (LLM) was used to assist in the writing and editing of this manuscript.
950 Specifically, we employed an LLM to help improve language expression, enhance readability, and
951 ensure clarity across all sections of the paper. The model provided support in tasks such as sentence
952 rephrasing, grammar checking, and improving overall textual fluency.

953 It is important to emphasize that the LLM did not contribute to the conception of research ideas,
954 methodology, or experimental design. All research concepts, insights, and analyses were indepen-
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