Grape Cold Hardiness Prediction via Multi-Task Learning

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Abstract

Cold temperatures during fall and spring have the potential to cause frost damage to grapevines and other fruit plants, which can significantly decrease harvest yields. To help prevent these losses, farmers deploy expensive frost mitigation measures such as sprinklers, heaters, and wind machines when they judge that damage may occur. This judgment, however, is challenging because the cold hardiness of plants changes throughout the dormancy period and it is difficult to directly measure. This has led scientists to develop cold hardiness prediction models that can be tuned to different grape cultivars based on laborious field measurement data. In this paper, we study whether deep-learning models can improve cold hardiness prediction for grapes based on data that has been collected over a 30-year time period. A key challenge is that the amount of data per cultivar is highly variable, with some cultivars having only a small amount. For this purpose, we investigate the use of multi-task learning to leverage data across cultivars in order to improve prediction performance for individual cultivars. We evaluate a number of multi-task learning approaches and show that the highest performing approach is able to significantly improve over learning for single cultivars and outperforms the current state-of-the-art scientific model for most cultivars.

Introduction

The ability of grapevines to survive cold temperatures during fall, winter, and spring, is known as Cold Hardiness (H_c) . Cold hardiness in grapes and other plants is dynamic in nature with a predictable seasonal trend. Cold hardiness is low at the beginning of fall as the plant has not yet acclimatized and peaks during mid-winter when the plant reaches acclimation. As spring arrives, the plant deacclimatizes and the cold hardiness decreases to the low summer levels. This means that during the fall and spring, when cold hardiness is low, unusually cold temperatures can be lethal, especially arising from sudden frost events.

To mitigate lethal damage due to cold temperatures, farmers can deploy expensive preemptive methods such as wind machines, sprinklers, and heaters to raise the air temperature. However, the decision of when to invest in expensive mitigation depends on knowledge of the current unknown

cold hardiness. While cold hardiness can be measured, it requires expertise and expensive equipment, which farmers rarely have. Thus, farmers rely on estimates of cold hardiness derived from a combination of experience and scientific models. This highlights the need for accurate data-centric models for cold-hardiness prediction.

Current state-of-the-art cold hardiness models (e.g. (Ferguson et al. 2014)) use a biological basis to obtain a parameterized model that can be tuned for different grape cultivars using cold-hardiness data. While reasonably effective, these models are relatively simple and only use ambient temperature as input. Rather, cold hardiness likely depends on multiple weather factors (e.g. humidity and precipitation) in complex ways (Mills, Ferguson, and Keller 2006) that are not full captured by current scientific models. This raises the question of whether modern machine learning methods can improve on current models via their increased expressiveness and ability to consume richer inputs.

In this work, we evaluate the use of Recurrent Neural Networks (RNNs) for predicting cold hardiness based on time series weather data. A key challenge is that groundtruth cold-hardiness data is quite limited in comparison with many applications of deep learning. In our experiments, we find that for some grape cultivars, where there is significant data, RNNs can be quite accurate and outperform a current state-of-the-art model. However, for cultivars with more limited data, the RNNs can perform poorly. This raises the question of whether we can leverage data across multiple cultivars to improve the prediction performance for cultivars with limited data.

Our main contributions are: 1) To frame this multi-cultivar learning problem as multi-task learning, and 2) To propose and evaluate a variety of multi-task RNN models on realworld data collected from over twenty cultivars with data amounts ranging from 34 to just 4 seasons. Our results show that multi-task learning is able to significantly outperform learning from just the data of a single cultivar and very often outperforms the state-of-the-art scientific model. We are aiming to install a model result from this work on an existing weather network for trial use by grape farmers.

Background

The cold hardiness of a plant characterizes its ability to resist injury during exposure to low temperatures. In this

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Figure 1: LTE_{50} measurements for three cultivars and temperature range for one season. The bowl shape indicates an initial increase in cold hardiness (decrease in lethal temperature) during acclimation followed by a decrease in cold hardiness during deacclimation. Notice the difference in variation of cold hardiness for the chosen cultivars.

work, we will focus on grapevine cold hardiness, where injury corresponds to lethal bud freezing, which decreases crop yield. In order to quantify how grapevine cold hardiness varies throughout the dormancy period, scientists use differential thermal analysis (DTA). DTA results in a measurement of the lethal temperatures at which 10%, 50%, and 90% of the bud population die/freeze, which are denoted by LTE_{10} , LTE_{50} , and LTE_{90} respectively (Mills, Ferguson, and Keller 2006). Figure 1 shows the LTE_{50} values for three cultivars and temperature ranges throughout a dormant season.

Since this measurement process requires expensive, specialized equipment and expertise, scientists have used collected data to develop grape cold-hardiness models, which aim to estimate the lethal temperatures based on only historical temperature data. The current state-of-the-art model, developed by Ferguson et al. (2011), integrates plant biology concepts to find a relation between daily temperatures and changes in cold hardiness. Intuitively, the Ferguson model computes the daily change in cold hardiness (e.g. as measured by LTE_{50}) based on the day's accumulated thermal time (being above or below certain temperature thresholds) weighted by coefficients that vary with the stage of dormancy. The model has a small number of parameters, e.g. thresholds for thermal times, which can be tuned for a particular cultivar. Tuning was done by performing a brute-force grid search over the parameter space to identify the parameter settings that resulted in the most accurate predictions. While this model has produced promising results and is in use by growers, it has limited expressiveness (only a handful of parameters) and only uses daily temperature data as input, rather than also factoring in other influential weather measurements (e.g. humidity and precipitation).

The limitations of the current scientific models raise the question of whether we can improve cold-hardiness prediction through the use of modern deep learning models. On one hand, such black-box deep models can be much more expressive and can easily incorporate additional weather data as input. On the other hand, the cold-hardiness data set sizes are relatively small from a deep-learning perspective, which may limit the potential benefits. The remainder of the paper explores this question. Below we first describe the cold-hardiness data sets used in our work followed by a description and evaluation of our deep learning approaches.

Cold Hardiness Datasets

The cold hardiness of endo–and ecodormant primary buds from up to 30 genetically diverse cultivars/genotypes of field-grown grapevines has been measured since 1988 in the laboratory of the WSU Irrigated Agriculture Research and Extension Center (IAREC) in Prosser, WA (46.29°N latitute; -119.74°W longitude). In the vineyards of the IAREC, the WSU-Roza Research Farm, Prosser, WA (46.25°N latitude; -119.73°W longitude), and in the cultivar collection of Ste. Michelle Wine Estates, Paterson, WA (45.96°N latitute; -119.61°W longitude), cane samples containing dormant buds were collected daily, weekly, or at 2-week intervals from leaf fall in autumn to bud swell in spring. These two phenological events typically occurred in October and in April, respectively (Ferguson et al. 2011, 2014).

All samples were analyzed with DTA to record ground truth for LTE_{10} , LTE_{50} , and LTE_{90} measurements of cold hardiness. Additionally, meteorological/environmental daily data from the closest on-site weather station to each vineyard (cultivar) was obtained using the API provided by AgWeatherNet (WSU 2022). The three stations used are Prosser.NE (46.25°N latitude; -119.74°W longitude), Roza.2 (46.25°N latitude; -119.73°W longitude), and Paterson.E (45.94°N latitude; -119.49°W longitude).

The result is a continually growing dataset for each cultivar that contains a varying number of seasons of daily weather data along with cold-hardiness LTE labels for the days that samples were collected. Following prior work we consider *a season* to extend from September 7th to May 15th, which is a conservative interval that should almost always contain the full dormancy period. Our experiments involve cultivars with data sets ranging from 34 to 4 seasons.

Cultivar Dataset Details. Table 1 shows a summary of the number of years of data collected for selected cultivars. The dataset for a given cultivar contains a row for each day of all data-collection seasons. Note, since cold hardiness was not measured on each day of a season, some rows do not contain LTE data. Below we highlight the key information contained in each row used by our models.

- DATE: The date of the weather observation.
- AWN_STATION: The closest AgWeatherNet station from where the environmental readings are taken.
- LTE values (when available): LTE_{10} , LTE_{50} , LTE_{90} . In degrees Celsius.
- MIN_AT, AVG_AT, MAX_AT: Minimum, average, and maximum air temperature observed at 1.5 meters above

Caltinga	LTE	LTE Total	LTE Total	
Cultivar	Data Seasons	Years of Data	Samples	
Barbera	2006-2022	14	151	
Cabernet Franc	2005-2012	4	35	
Cabernet Sauvignon	1988-2022	34	829	
Chardonnay	1996-2022	26	783	
Chenin Blanc	1988-2022	18	193	
Concord	1988-2022	27	484	
Gewurztraminer	2005-2016	9	101	
Grenache	2006-2022	14	151	
Lemberger	2006-2016	6	60	
Malbec	2004-2022	17	261	
Merlot	1996-2022	26	897	
Mourvedre	2005-2022	12	133	
Nebbiolo	2006-2022	14	152	
Pinot Gris	2003-2022	17	190	
Riesling	1988-2022	34	636	
Sangiovese	2005-2022	15	165	
Sauvignon Blanc	2006-2022	12	140	
Semillon	2006-2022	13	201	
Syrah	1999-2022	23	486	
Viognier	1999-2022	18	206	
Zinfandel	2006-2022	14	150	

Table 1: Summary of cultivars' LTE data collection.

the ground. In degrees Celsius.

- MEAN_AT: $(MIN_AT + MAX_AT)/2$. In degrees Celsius.¹
- MIN_RH, AVG_RH, MAX_RH: Minimum, average, and maximum relative humidity value observed at 1.5 meters above the ground. In percent.
- MIN_DEWPT, AVG_DEWPT, MAX_DEWPT: Minimum, average, and maximum dew point (temperature the air needs to be cooled to in order to achieve relative humidity). In degrees Celsius.
- P_INCHES: Observed sum of precipitation for the daily period. In inches.
- WS_MPH, MAX_WS_MPH: Average and maximum observed wind speed at 1.5 meters above the ground for the daily period. In Miles Per Hour.

Deep Cold-Hardiness Models and Training

Given the availability of cold-hardiness data, we can formulate cold-hardiness prediction as a sequence prediction problem. We will use *i* to index the different grape cultivars with N_i denoting the number of seasons available for cultivar *i*. The sequence data for season *k* of cultivar *i* is denoted by $S_{i,k}$ and has the form $S_{i,k} = (x_1, y_1, x_2, y_2, \ldots, x_H, y_H)$, where x_t is the weather data for day *t*, y_t is the ground truth LTE data for day *t*, and *H* is the number of days per season. Recall that y_t is not measured on each day of a season (e.g. measured every two weeks) and hence for days where the LTE measurements are unavailable $y_t = N/A$. Finally, the data set for cultivar *i* is denoted by $D_i = \{S_{i,k} \mid k \in \{1, \ldots, N_i\}\}$.

Given a data set D_i our learning goal is to produce a model M_i that can take as input a sequence of daily weather measurements (x_1, x_2, \ldots, x_t) up to a particular day t and produce a sequence of predicted LTE estimates $(\hat{y}_1, \hat{y}_2, \ldots, \hat{y}_t)$ for cultivar i. Typically, a farm manager will be most interested in the estimate \hat{y}_t . This estimate can then be compared to the low-temperature forecast for that day to help decide whether to prepare for frost mitigation measures. The key question of this work is to evaluate whether modern deep learning methods can provide farm managers with improved predictions compared to the current state-of-the-art cold-hardiness models.

We will refer to the problem of learning M_i based on only D_i as single-task learning (STL), which is the general framework used for the vast majority of deep learning applications. Importantly, the performance of STL is significantly influenced by the amount of available training data, which according to Table 1 varies widely across the different cultivars. Thus, we might expect STL performance for cultivars with small datasets to suffer in comparison to those with large datasets. To address this issue, we consider the multi-task learning (MTL) framework (Caruana 1997; Vandenhende et al. 2021; Crawshaw 2020; Zhang and Yang 2021), which involves learning a predictive model for cultivar i using a combined dataset of all cultivars $D = \{D_1, D_2, \dots, D_C\}$, where C is the number of cultivars for which we have data. Intuitively, MTL offers the potential to identify common structures among the multiple learning tasks (i.e. cultivars) in order to improve performance for individual cultivars, especially those with limited data.

Below, we first introduce the STL deep model that we developed, which will serve as our deep-learning baseline for cold-hardiness prediction. Next, we introduce two frameworks for modifying that model to support MTL. Finally, we describe certain details of the training strategy used in our experiments. To the best of our knowledge, this is the first work that has considered deep models for cold-hardiness prediction in both the STL and MTL settings.

Single-Task Model

Our basic STL makes causal LTE predictions by sequentially processing a weather data sequence x_1, x_2, \ldots, x_t and at each step outputting the corresponding LTE estimate. For this purpose we use a recurrent neural network (RNN) model (Rumelhart, Hinton, and Williams 1985), which is a widely used model for sequence data. The RNN backbone used by both our STL and MTL models is illustrated in Figure 2a), which we denote by f_{θ} with parameters θ . The backbone network begins with two Fully Connected (FC) layers, followed by a Gated Recurrent Unit (GRU) layer (Cho et al. 2014), which is followed by another FC layer. Our STL model, shown in Figure 2b), simply feeds daily weather data x_t into the first FC layer as input and adds an additional FC layer to produce the final LTE prediction output. Intuitively, the GRU unit, through its recurrent connection is able to build a latent-state representation of the sequence data that has been processed so far. For our cold-hardiness problem, this representation should capture information about the weather history which is useful for predicting LTE. In some sense, the latent state can be thought of as implicitly approximating the internal state of the plant as it evolves during dormancy. As described below, each STL model M_i

¹This is recorded since it is the temperature measure used by the scientific model.



Figure 2: Network Architectures. FC denotes fully connected layers and GRU denotes Gated Recurrent Unit. a) The RNN backbone is used to process weather data sequences (x_t) . b) The single-task model with a single prediction head for a single cultivar. c) Multi-Head Model which has a prediction head for each cultivar allowing backbone features to be shared. d) Task Embedding Model, which combines the weather data features with a learned task embedding for each cultivar before entering the backbone network.

is trained independently on its cultivar-specific dataset D_i .

Multi-Task Models

We consider two types of MTL models that directly extend the RNN backbone of Figure 2a), the multi-head model and the task embedding model.

Multi-Head Model. The multi-head model is perhaps the most straightforward approach to MTL and has been quite successful in prior work when tasks are highly related (Caruana 1997). As illustrated in Figure 2c), the multi-head model is identical to the STL model, except, that it adds C parallel cultivar-specific fully-connected layers to the backbone (i.e. prediction heads). Each prediction head is responsible for producing the LTE prediction for its designated cultivar. This model allows the cultivars to share the features produced by the RNN backbone, with each cultivar-specific output simply being a linear combination of the shared features. Intuitively, if there are common underlying features that are useful across cultivars, then this architecture allows those to emerge based on the combined set of data. Thus, cultivars with small amounts of data can leverage those useful features and simply need to tune a set of linear weights based on the available data. We abbreviate this model with MultiH in future sections.

Task Embedding Models. Our proposed Task Embedding models are motivated by thinking about current scientific models and how they address multiple tasks. The Ferguson model, for example, has a fixed structure, based on scientific knowledge, but a small number of parameters that can be tuned for each cultivar. Our task embedding model aims to generalize this concept by having a neural network learn both the structure of the model that accepts task-specific parameters as well as learning the parameters of each cultivar. Note that the cultivar parameters and model structure will not have a clear scientific interpretation due to the black-box nature of deep models. The trade-off for interpretability is the potential for better performance due to increased expressive power.

Specifically, our proposed Task Embedding models, as shown in Figure 2d), are similar in spirit to context-sensitive

neural networks (Silver, Poirier, and Currie 2008; Schreiber, Vogt, and Sick 2021; Schreiber and Sick 2021), where a task-specific context is provided as additional input to the neural network with only a single output being computed.

We obtain this task-specific context by encoding the task as a one-hot vector and finding a corresponding mapping using a differentiable embedding layer. We explore different ways of incorporating the obtained task embedding, via element-wise Addition, Concatenation, and element-wise Multiplication. We abbreviate these models with AddE, ConcatE, MultE in future sections.

Model and Training Details

We construct our dataset by selecting the dormant season data for all cultivars. Missing features are filled in by linear interpolation. We discard seasons that have < 10% valid LTE readings. We only include seasons where at least 90% of temperature data is not missing. Missing LTE label readings are not interpolated, instead, the missing LTE labels' losses are masked during the training and evaluation process. We choose 2 seasons for each cultivar as our test set. We run three trials of training for all our experiments with different train/test splits and average the performance over the three trials.

We rely on the following weather features for learning our models - Temperature, Humidity, Dew Point, Precipitation, and Wind Speed. Our models output predictions for LTE_{10} , LTE_{50} , and LTE_{90} which are optimized simultaneously, helping in inductive transfer. We focus exclusively on the model's predictive power for LTE_{50} in this work. We consider the Mean Squared Error(MSE) as our loss function and treat the Root Mean Squared Error(RMSE) as our performance metric. We consider Adam (Kingma and Ba 2014) as the optimizer of choice for our training process. We use a learning rate of 0.001 with a batch size of 12 seasons shuffled randomly. We train all our models for 400 epochs. The input features have a dimensionality of 12. The output dimensionality of the linear layers of the RNN backbone are 1024, 2048 and 1024 respectively. The GRU has a hidden state and internal memory of dimensionality 2048.

Cultivar	MultE	ConcatE	AddE	MultiH	Single	Ferguson
Barbera	1.92	1.50	2.07	1.89	4.22	1.78
Cabernet Franc	4.84	2.36	3.49	2.39	4.00	1.45
Cabernet Sauvignon	2.93	1.75	1.82	2.27	3.43	1.83
Chardonnay	1.33	1.46	1.44	1.40	1.60	1.79
Chenin Blanc	1.85	1.51	1.57	1.45	2.47	2.27
Concord	2.33	2.42	2.32	1.98	2.61	2.02
Gewurztraminer	1.97	1.40	1.66	1.20	2.70	1.84
Grenache	3.07	1.86	2.17	1.79	2.86	1.92
Lemberger	3.01	1.65	2.24	1.49	3.23	2.21
Malbec	1.80	1.32	1.32	0.96	1.71	1.66
Merlot	1.74	1.53	1.39	1.53	1.66	1.55
Mourvedre	1.84	1.65	1.70	1.56	2.25	1.83
Nebbiolo	2.36	1.58	1.87	1.24	2.48	1.80
Pinot Gris	2.07	1.61	1.63	1.61	2.04	2.02
Riesling	2.80	1.47	1.77	1.97	3.63	1.55
Sangiovese	1.65	1.73	1.71	1.40	1.84	1.61
Sauvignon Blanc	1.33	1.43	1.52	1.22	1.71	1.42
Semillon	2.37	1.67	1.42	1.75	3.58	1.50
Syrah	1.22	1.22	1.28	1.29	1.57	1.25
Viognier	3.90	1.75	2.30	2.28	4.16	1.36
Zinfandel	3.10	1.45	1.56	1.60	2.64	1.90

Table 2: Comparison of the performance of proposed MTL methods with STL and the existing state-of-the-art method. Note that the performance is measured in terms of Root Mean Squared Error.

Experiments

In this section, we present our main empirical results. Our experiments involve 21 cultivars from Table 1. In particular, we removed any cultivar that has less than 4 years of data and removed cultivars for which the Ferguson model results were unavailable for comparison.

Multi-Task Versus Single-Task Learning. Table 2 shows the root mean squared error (RMSE) of the multi-task models, single-task model, and Ferguson model for all 21 cultivars. The first observation is that with the exception of Chardonnay, the single-task model never outperforms the state-of-the-art Ferguson model. For cultivars with small amounts of data, there is often a dramatic decrease in performance over Ferguson, while other cultivars with larger datasets are close to Ferguson's performance.

The second observation is that for each cultivar, with rare exceptions, the 4 multi-task models all outperform the corresponding single-task model. As expected, the improvement tends to be most pronounced for the smaller dataset cultivars. This shows that multitask learning is indeed able to identify and exploit common structures among the different cultivars, leading to improved generalization. Among the MTL methods, the MultiHead approach consistently outperforms other methods. Among the task embedding approaches, the concatenation approach performs best.

We observe, with the exception of Cabernet Franc and Viognier, that our approaches outperform the Ferguson model, the MultiHead or concat task embedding approaches being the best-performing models for most cultivars. The gap in the performance is more dramatic in cultivars with low data, such as Lemberger and Gewurztraminer.

Impact of Task Dataset Size. Table 3 shows the performance of MTL(MultiHead) and STL models when we choose one of three cultivars (Riesling, Merlot, Cabernet Sauvignon) with \sim 30 seasons of data and artificially select

Cultivar	2	5	10	20	All
Riesling (MTL)	2.25	2.16	2.10	1.71	1.97
Riesling (STL)	4.59	4.40	3.66	3.41	3.63
Cabernet Sauvignon (MTL)	2.02	2.16	2.27	2.07	2.27
Cabernet Sauvignon (STL)	2.68	3.24	3.68	2.91	3.43
Merlot (MTL)	1.69	1.54	1.55	1.41	1.53
Merlot (STL)	2.26	1.99	1.83	1.67	1.66

Table 3: Measuring the impact of varying the dataset size for chosen cultivars. The experiment is conducted for both STL and MTL. We choose 2, 5, 10, and 20 seasons as reasonable choices to evaluate. The performance is measured in terms of RMSE.

only a subset of seasons, and train the MultiHead model. We also train STL models in the same setup. As expected for STL, with the exception of Cabernet Sauvignon, introducing more seasons of data up to an extent does help in improving performance.²

Interestingly we see that an MTL model trained on just 2 or 5 seasons of data for a cultivar outperforms an STL model using all of that cultivar's data. This reflects the fact that MTL is indeed able to leverage the information present in other cultivars to learn a good model for that specific cultivar. In a sense, the data from other cultivars appear to be as valuable as tens of seasons of data for a specific cultivar.

Impact of the number of tasks - The goal here is to understand how different subsets of tasks impact the performance of an MTL model. Here, we select different subsets of our tasks, 10 tasks with the most amount of data, 10 tasks with the least amount of data, and 10 tasks with a mix of

²Note that there is a consistent decrease in performance when going from 20 seasons to ALL. The reasons for this remain to be explored; however, it is likely due to the influence of a small number of unusual seasons.

Cultivar	High	Low	Mix	All	Single
Riesling	1.95		1.70	1.97	3.63
Cabernet Sauvignon	2.49		2.28	2.27	3.43
Chardonnay	1.33		1.41	1.40	1.60
Concord	2.00		1.90	1.98	2.61
Merlot	1.46		1.39	1.53	1.66
Syrah	1.13			1.29	1.57
Chenin Blanc	1.60			1.45	2.47
Viognier	2.60			2.28	4.16
Malbec	1.12			0.96	1.71
Pinot Gris	1.48			1.61	2.04
Barbera		2.93		1.89	4.22
Grenache		2.04		1.79	2.86
Nebbiolo		1.85		1.24	2.48
Zinfandel		2.07		1.60	2.64
Semillon		2.47		1.75	3.58
Mourvedre		1.93	1.75	1.56	2.25
Sauvignon Blanc		1.65	1.27	1.22	1.71
Gewurztraminer		1.83	1.35	1.20	2.70
Lemberger		2.24	1.50	1.49	3.23
Cabernet Franc		2.91	2.13	2.39	4.00

Table 4: Measuring the impact of choosing a subset of available tasks for MTL and how it fares against choosing all tasks.

high and low amounts of data. We train the MultiHead architecture for this experiment.

Table 4 presents the cultivars in order of largest to smallest datasets. Each column corresponds to the subset of cultivars used in each experiment. Interestingly, we observe that an MTL model trained on any of our chosen subsets always outperforms single-task models for all cultivars.

For cultivars with relatively higher amounts of data, surprisingly, it is better to choose a mix of high and low data cultivars to get a better performing model. Including all the cultivars for training does not lead to consistent gains for all cultivars with high amounts of data. The reasons for this observation require further analysis and experimentation.

For cultivars with lower amounts of data, again, choosing a mix of high and low data cultivars leads to a better performing model. Including all the cultivars for training does indeed lead to consistent gains over choosing a subset.

Impact of Training Setting. Here, we consider a different training setting, Transfer Learning (Bozinovski and Fulgosi 1976), where a new cultivar arrives and is incorporated into the model without access to past data. This is in contrast to MTL where all datasets are accessible at training time. We consider finetuning as a straightforward approach to transfer learning. Finetuning, in the case of MultiHead refers to replacing the task-specific final layers with a newly initialized layer for the new task. For the task embedding approaches, finetuning refers to learning the coefficients of a linear combination of existing task embeddings.

Table 5 shows the RMSE metrics for the finetuning paradigm for the different proposed methods relative to their corresponding RMSE metrics for MTL from Table 2.

In the case of the MultiHead architecture, we observe that finetuning is on par with MTL. This seems to indicate that there are no tasks that hurt the MTL training process.

Although, for the Task Embedding approaches, we see

Cultivar	ConcatE FT	MultE FT	AddE FT	MultiH FT
Barbera	-1.02	0.04	-1.69	0.01
Cabernet Franc	-1.50	2.41	-2.28	-0.05
Cabernet Sauvignon	-1.02	0.64	-1.29	-0.01
Chardonnay	-1.14	0.04	-3.01	0.11
Chenin Blanc	-0.74	0.35	-2.96	-0.05
Concord	-3.38	0.11	-2.96	-0.24
Gewurztraminer	-1.58	0.51	-2.66	-0.25
Grenache	-0.32	1.27	-1.69	-0.01
Lemberger	-2.53	1.37	-0.43	-0.16
Malbec	-2.78	0.77	-1.66	-0.07
Merlot	-0.84	0.32	-2.32	0.11
Mourvedre	-1.28	0.22	-1.58	-0.07
Nebbiolo	-2.51	0.72	-1.31	-0.41
Pinot Gris	-1.14	0.53	-2.44	0.08
Riesling	-1.78	1.14	-1.00	0.31
Sangiovese	-0.96	0.31	-1.73	0.05
Sauvignon Blanc	-1.22	0.09	-0.23	-0.02
Semillon	-1.46	0.97	-3.08	0.35
Syrah	-0.97	-0.06	-1.87	0.00
Viognier	-1.25	2.13	-1.89	0.52
Zinfandel	-2.24	1.31	-1.75	-0.19
Median	-1.25	0.53	-1.75	-0.01
Mean	-1.51	0.72	-1.90	0.00

Table 5: Comparing Transfer Learning with Multi-Task Learning. Note that the performance is relative to corresponding MTL counterparts in table 2. If a term is positive, it means that transfer learning does better than MTL in that case. We abbreviate finetuning with FT in the column names.

that finetuning does worse than MTL for most cultivars for the Concatenate and Additive variants. For the multiplicative embedding variants, we see marginal to substantial gains in performance.

Path to Deployment

Farmers use AgWeatherNet (WSU 2022) and WSU Viticulture and Enology (Mills 2022) websites to monitor cold hardiness through the deployment of the Ferguson model and publication of real LTE values, respectively. Our goal is to finalize the MTL-based models proposed in this paper and deploy them onto AgWeatherNet for the 2022-2023 season for beta testing.

Conclusion

We showed that multi-task learning is an effective approach to predicting Cold Hardiness for grapevines. In particular, our model consistently outperforms the state-of-the-art scientific model without relying on expert domain knowledge. This model will be deployed on an existing weather network for the 2022-2023 season. In the future, we plan to apply these ideas to cold-hardiness prediction for other crops, such as cherries and apples. In addition, we plan to investigate the utility of MTL for other agriculture-related problems with limited data.

Acknowledgements

This research was supported by USDA NIFA award No. 2021-67021-35344 (AgAID AI Institute). The authors thank Lynn Mills and Alan Kawakami for the LTE data collection.

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