Money Doesn’t Grow on Trees, But Forecasts Do: Forecasting Extreme Precipitation with Random Forests

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ABSTRACT
Approximately eleven years of reforecasts from NOAA’s Second Generation Global Ensemble Forecast System Reforecast (GEFS/R) model are used to train a contiguous United States (CONUS)-wide gridded probabilistic prediction system for locally extreme precipitation, developed primarily using the random forest (RF) algorithm. Locally extreme precipitation is quantified for 24-hour precipitation accumulations in the framework of average recurrence intervals (ARIs), with two severity levels: 1- and 10-year ARI exceedances. Forecasts are made from 0000 UTC forecast initializations for two 1200 UTC–1200 UTC periods: Days 2 and 3 comprising respectively forecast hours 36–60 and 60–84. Separate models are trained for each of eight forecast regions and for each forecast lead time. GEFS/R predictors vary in space and time relative to the forecast point, and include not only the quantitative precipitation forecast (QPF) output from the model, but also variables that characterize the meteorological regime, including winds, moisture, and instability. Numerous sensitivity experiments are performed to determine the effects of the inclusion or exclusion of different aspects of forecast information in the model predictors, the choice of statistical algorithm, and the effect of performing dimensionality reduction via principal component analysis as a pre-processing step. Overall, it is found that the machine learning (ML)-based forecasts add significant skill over exceedance forecasts produced from both the raw GEFS/R ensemble QPFs and from the European Centre for Medium-Range Weather Forecasts’ (ECMWF) global ensemble across almost all regions of CONUS. ML-based forecasts are found to be underconfident, while raw ensemble forecasts are highly overconfident.
1. Introduction

Locally extreme precipitation can cause a variety of costly, disruptive, and endangering impacts, including flooding, flash flooding, and landslides. In 2016 alone, these hazards combined caused more than 120 fatalities and $10 billion in damages over the United States (NWS 2017b). The prediction of flash floods is a notoriously challenging forecast problem, requiring not only accurate prediction of heavy rainfall magnitudes, but also of the spatiotemporal distribution of that rainfall; the hydrologic interactions between precipitation, terrain, and the land surface; and also of antecedent precipitation and its effects on soil conditions. Forecasting precipitation processes responsible for most observed extreme rainfall over the contiguous United States (CONUS) is often considered among the most challenging problems in contemporary numerical weather prediction (NWP; e.g. Fritsch and Carbone 2004; Novak et al. 2014). Given that the rainfall forecast alone presents such a considerable challenge, the additional hydrologic considerations in the flash flood forecast problem present an even more daunting task. While recent advances in heavy rainfall and flash flood forecasting have been made (e.g. Hapuarachchi et al. 2011; Novak et al. 2014; Barthold et al. 2015), forecasts still struggle in many situations (e.g. Delrieu et al. 2005; Lackmann 2013; Schumacher et al. 2013; Gochis et al. 2015; Nielsen and Schumacher 2016, among many others) and substantial progress remains to be made.

Contemporary operational dynamical forecast models often struggle to simulate accurately the physical processes responsible for extreme precipitation production. For example, models with parameterized convection often have a variety of persistent errors and biases associated with their depiction of convective systems, which are responsible for the majority of flooding rains over much of CONUS (e.g. Schumacher and Johnson 2006; Stevenson and Schumacher 2014; Herman and Schumacher 2016a). These include a tendency to underpredict total rainfall from convective
systems (e.g. Schumacher and Johnson 2008; Herman and Schumacher 2016a); produce systems
displaced too far to the north and west from where they are observed (e.g. Grams et al. 2006;
Wang et al. 2009; Clark et al. 2010); initiate convection too early (e.g. Davis et al. 2003; Wilson
and Roberts 2006; Clark et al. 2007); generate systems with too large an areal extent (e.g. Wilson
and Roberts 2006); and propagate them incorrectly, too slowly, or not at all (e.g. Davis et al. 2003;
Pinto et al. 2015). While convection-allowing models (CAMs) can better resolve the physical
processes responsible for heavy rainfall generation (e.g. Kain et al. 2006; Weisman et al. 2008;
Duda and Gallus 2013), they too can suffer from many of these biases (e.g. Kain et al. 2006; Lean
et al. 2008; Kain et al. 2008; Weisman et al. 2008; Herman and Schumacher 2016a). Furthermore,
although there is a plethora of CAM guidance out to the day-ahead time frame (out to 36 hours
to perhaps 48 hours after initialization), due to current computational constraints, there is almost
no operational CAM guidance running out to two days ahead, and nothing operational that runs
to three days ahead or beyond. Instead, global ensembles with parameterized convection serve
as the primary source of forecast information and uncertainty quantification at these lead times.
Nevertheless, there is considerable utility in skillful extreme precipitation forecasts at these longer
lead times, since many mitigative actions that may not be feasible to execute in a matter of hours,
but easily accomplished with a day or more of warning. Statistical post-processing of global
ensemble output can potentially alleviate many of these dynamical model deficiencies and provide
skillful extreme precipitation guidance at medium-range time scales. A specific focus on the Day
2–3 period is warranted due to the increased existing operational emphasis on these lead times
compared with even longer ones, such as the Excessive Rainfall Outlooks produced by the Weather
Prediction Center (Barthold et al. 2015) which forecast locally excessive rainfall across CONUS
for Days 1–3.
There is a long history of successful application of statistical post-processing to dynamical model output (e.g. Klein et al. 1959; Glahn and Lowry 1972). Model Output Statistics (MOS; e.g. Glahn and Lowry 1972), is a simple, effective multivariate linear regression technique relating a set of dynamical model predictors to sensible weather predictands such as minimum and maximum temperature, wind speeds, and precipitation probability. This basic technique has long demonstrated skill over both the underlying models and even human forecasters (e.g. Jacks et al. 1990; Vislocky and Fritsch 1997; Hamill et al. 2004; Baars and Mass 2005), but is inherently limited by the linear assumptions underlying the method. Statistical post-processing techniques have also been successfully applied to QPFs, from early linear approaches (e.g. Bermowitz 1975; Antolik 2000) to more contemporary techniques that can exploit more complex variable relationships, including neural networks (e.g. Hall et al. 1999), reforecast analogs (e.g. Hamill and Whitaker 2006; Hamill et al. 2015), logistic regression (LR; e.g. Applequist et al. 2002; Whan and Schmeits 2018), random forests (RF; e.g. Gagne et al. 2014; Ahijevych et al. 2016; Gagne et al. 2017; Whan and Schmeits 2018), and other parametric techniques (e.g. Scheuerer and Hamill 2015; Whan and Schmeits 2018). For other meteorological applications, other machine learning algorithms, such as support vector machines (e.g. Zeng and Qiao 2011; Herman and Schumacher 2016b) and boosting (e.g. Herman and Schumacher 2016b; Hong et al. 2016) have also been successfully applied. Related techniques have also been applied to forecasting related high-impact phenomena, such as severe hail (Brimelow et al. 2006; Gagne et al. 2015) and tornadoes (Alvarez 2014). One of the most powerful aspects of machine learning algorithms—and RFs in particular—is finding patterns and non-linear interactions in the supplied training data (e.g. Breiman 2001). Depending on the extent and diversity of the data supplied in these experiments, trained RFs pose the theoretical capability of diagnosing and automatically correcting for various kinds of model biases, including context-dependent quantitative biases, such as QPF being systematically too high or too low; spa-
tial displacement biases in the placement of extreme precipitation features; and, to some extent, temporal biases in the initiation or progression of extreme precipitation features.

This study makes a comprehensive investigation of using a global reforecast dataset to produce skillful and reliable probabilistic forecasts of locally extreme precipitation using the RF statistical post-processing technique in the medium-range. The following section provides further background and rigorously describes the data and methods used, algorithms employed, models trained, and experiments performed. Section 3 presents results of the sensitivity experiments conducted, while Section 4 presents the final results of the trained models and provides two brief case studies illustrating the process. Section 5 summarizes the findings of this study, outlines complementary analysis of these models, identifies avenues for further research, and discusses the broader implications of the results on numerical weather prediction and post-processing.

2. Data and Methods

There are several successive steps applied in creating the final forecasts evaluated in this study. A schematic overview of the forecast pipeline for the models trained in this study is depicted in Figure 1. Many types of hydrometeorological information are first taken, then assembled in a methodical manner, further pre-processed for subsequent analysis, analyzed using a statistical machine learning algorithm, and finally, extreme precipitation forecast guidance is produced and evaluated. This section details each of these steps in the model development and evaluation process.

a. Datasets

Dynamical model data used for training the RF models in this study comes from NOAA’s Second-Generation Global Ensemble Forecast System Reforecast (GEFS/R; Hamill et al. 2013)
dataset. The GEFS/R is a global 11-member ensemble with parameterized convection and T254L42 resolution—which corresponds to an effective horizontal grid spacing of ≈55 km at 40° latitude—initialized once daily at 0000 UTC back to December 1984. Perturbations are applied only to the initial conditions, and are made using the ensemble transform with rescaling technique (Wei et al. 2008). The ensemble system used to generate these reforecasts is nearly static throughout its 30+ year period of coverage, though updates to the operational data assimilation system over time have resulted in some changes in the bias characteristics of its forecasts over the period of record (Hamill 2017). Some forecast fields are preserved on the native Gaussian grid (≈0.5° spacing), while others are available only on a 1° × 1° grid. Temporally, forecast fields are archived every three hours out to 72 hours past initialization, and are available every six hours beyond that.

This study employs an almost 11-year period of record to explore this forecast problem, using daily initializations from January 2003 through August 2013.

In creating probabilistic extreme precipitation forecast guidance, the predictand must first be concretely specified and a robust, consistent verification framework established. One of the many challenges in heavy rainfall and flash flood forecasting is the considerable difficulty in verifying events (e.g. Welles et al. 2007; Gourley et al. 2012; Barthold et al. 2015), as every approach has its deficiencies and limitations. It is attractive to consider the problem from a simple perspective of quantitative precipitation estimate (QPE) exceedances of some temporally static threshold. In particular, a fixed threshold (e.g. 50 mm hr$^{-1}$) can be used as a proxy for flash flooding (e.g. Brooks and Stensrud 2000; Hitchens et al. 2013), as can exceedances of thresholds defined relative to the local precipitation climatology (e.g. Schumacher and Johnson 2006; Stevenson and Schumacher 2014; Herman and Schumacher 2016a), such as average recurrence intervals (ARIs). An ARI defines a fixed frequency relative to the hydrometeorological climatology of the region; in particular, it corresponds to the expected duration, given the local climatology, between exceedances of a
given threshold. For example, the 1-year ARI for 24-hour precipitation accumulations describes the accumulation amount for which one would expect the mean duration between exceedances of said amount to be one year. Past research has shown that a fixed-frequency ARI-based framework has better correspondence with heavy precipitation impacts than the use of any fixed threshold across the hydrometeorologically diverse regions of CONUS (e.g. Reed et al. 2007). From the perspective of forecast verification, defining extreme precipitation with respect to a fixed threshold exceedance raises challenges when applied uniformly across CONUS. For example, skill differences observed between regions may simply be an artifact of a regionally varying event climatology rather than “true” regional differences in forecast skill (e.g. Hamill and Juras 2006). The ARI framework avoids this issue and provides reasonable correspondence with precipitation impacts while avoiding the additional complications such as antecedent conditions, local hydrology, and urban effects (e.g. Herman and Schumacher 2016a) and is consequently used to quantify extreme rainfall for this study.

Specifically, forecast probabilities are issued for 24-hour ARI exceedances at each GEFS/R archive grid point on its native Gaussian grid at all points across CONUS, using a predictand with three categories: 1) No 1-year ARI exceedance at any point within the grid point domain, 2) At least one 1-year ARI exceedance, but no 10-year ARI exceedances within the grid point domain, and 3) At least one 10-year ARI exceedance within the grid point domain. For evaluation, probabilities from the middle and most severe categories are often aggregated to produce a 1-year ARI exceedance probability. This approach has the advantage of retaining aspects of the anticipated event severity as would be retained in a regression context but is largely lost when performing single category classification. While there can be some additional complications especially with respect to calibration, formulating the prediction problem as a single multicategory classification task rather than multiple distinct binary category models also ensures mathematical consistency.
of the exceedance probabilities within the generated probability mass functions in a way that the latter approach would not.

In aggregating multiple QPE-to-ARI threshold grid point comparisons in a single predictand, the forecasts issued correspond to neighborhood event probabilities, an increasingly popular method of communicating probabilistic high-impact weather information in forecast operations (e.g. Barthold et al. 2015; NWS 2017a). Counting any one of several possible point exceedances as an “event” results in the event having a higher observed relative frequency relative to that of any of the individual point exceedances; the event frequency in this framework thus exceeds the purported frequency suggested by the ARI. However, the fixed-frequency property, and thus many of the aforementioned desirable properties of the framework, are approximately retained. For this study, focus is placed exclusively on two 24-hour forecast periods: the 1200–1200 UTC period corresponding to forecast hours 36–60 from the GEFS/R forecast fields and the subsequent 24-hour period encompassing forecast hours 60–84, denoted respectively as Days 2 and 3. At these times, there is typically some knowledge to characterize the environmental conditions in which precipitation may form, but it is beyond the current range of operational CAM guidance.

Verification comes from the National Centers for Environmental Prediction (NCEP) Stage IV Precipitation Analysis (Lin and Mitchell 2005) QPE product, created operationally since December 2001. Stage IV provides 24-hour analyses over the CONUS on a ∼4.75 km grid. It uses both rain gauge observations and radar-derived rainfall estimates to generate an analysis, and is further quality controlled via NWS River Forecast Centers (RFCs) to ensure stray radar artifacts and other spurious anomalies do not appear in the final product. Despite some limitations (Herman and Schumacher 2016a; Nelson et al. 2016), its analysis quality; resolution, allowing better ability to capture precipitation extremes compared with other QPE products (e.g. Hou et al. 2014); and data record length make it preferable to analogous products.
The ARI thresholds associated with the 1- and 10-year ARIs for 24-hour precipitation accumulations are generated using the same methodology as Herman and Schumacher (2016a), where CONUS-wide thresholds are produced by stitching thresholds from several sources. NOAA’s Atlas 14 thresholds (Bonnin et al. 2004, 2006; Perica et al. 2011, 2013), an update from older work and currently under development, are used wherever they were available at the commencement of this study. For five northwestern states—Washington, Oregon, Idaho, Montana, and Wyoming—updated thresholds are not available, and derived NOAA Atlas 2 threshold estimates are used instead (Miller et al. 1973). Additionally, in Texas and the Northeast—New York, Vermont, New Hampshire, Maine, Massachusetts, Connecticut, and Rhode Island—Technical Paper 40 (TP-40; Hershfield 1961) thresholds are used; everywhere else uses the Atlas 14 threshold estimates. The 10-year ARI thresholds (Fig. 2b) show a similar spatial pattern to the 1-year ARI thresholds (Fig. 2a), but are substantially higher everywhere. More significantly, it is apparent that at both severity levels, there are large regional disparities in the threshold magnitudes. Over climatologically wet regions of CONUS, such as the Pacific coastal mountains and immediately along the Gulf Coast, thresholds are as high as 100–150 mm and 250–300 mm for 1-year and 10-year ARIs, respectively. Over central and eastern CONUS, thresholds tend to decrease smoothly with increasing latitude and distance from major bodies of water. Sharper variations are seen in areas of complex terrain over western CONUS. In the driest parts of the Arid Southwest and Intermountain West, thresholds can be as low as 10–15 mm and 25–30 mm for the two ARI levels—a full order of magnitude difference from the largest thresholds at the same intensity level.

Forecast models in this study are trained separately for eight distinct, yet cohesive and internally fairly hydrometeorologically homogeneous regions of CONUS, using the delineation indicated

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1The northeastern states did receive updated Atlas 14 estimates in October 2015, but TP-40 thresholds were retained for consistency with prior work.
in Figure 3. Observed 1- and 10-year ARI exceedance events that occurred during the period of
record (Fig. 2c,d) highlight important regional differences in the seasonal climatology of ARI
exceedances across CONUS. In the Pacific Coast (PCST) region, the vast majority of exceedances
at both the 1-year and 10-year severity levels occur in the cool-season, and occur largely from
atmospheric river events with large moisture transport impinging on coastal topography (e.g. Rutz
et al. 2014; Herman and Schumacher 2016a). This seasonality holds to a lesser extent in the
neighboring Southwest (SW) region, with some signal carrying over to the Rockies (ROCK) region
as well. In the central and eastern regions, the majority of events occur during the warm-season
from more scattered convective-scale processes, particularly in the months of May, June, and July
(e.g. Schumacher and Johnson 2006; Herman and Schumacher 2016a). Tropical cyclones can
cause widespread and very significant rainfall, and comprise a substantial portion of the extreme
precipitation climatology, especially in the Northeast (NE) and Southeast (SE) regions. Due to
the spatial extent of their impacts and immense rainfall totals they can produce, they form a much
larger fraction of the climatology of 10-year ARI exceedances (Fig. 3d) than 1-year events (Fig.
3c). Additionally, the numbers are lower than would be expected; by the explicit exceedance
frequencies associated with the thresholds, one would expect an average of one exceedance per
point per year over the period of record for the 1-year events (Fig. 3c) and 0.1 exceedances for 10-
year events (Fig. 3d); in reality, event counts are only approximately half of that. This is consistent
with previous findings (e.g. Herman and Schumacher 2016a), and likely in part attributable to
limitations in the Stage IV product to capture extremes (e.g. Nelson et al. 2016). There is also
quite a bit of region-to-region variability in event counts, particularly for 10-year exceedances,
much of which is attributable to statistical variability from having a short data record in relation to
the event frequency.
b. Predictor Assembly

Input predictors, or features, to the random forests can be partitioned into two categories: model predictors and background predictors; the former constitute the vast majority of inputs. Model predictors come from atmospheric fields forecast in the GEFS/R which bear a known physical relationship with extreme precipitation. A core set of \( f = 9 \) fields used in this study are: accumulated precipitation (APCP), convective available potential energy (CAPE), convective inhibition (CIN), precipitable water (PWAT), surface temperature (T2M) and specific humidity (Q2M), surface zonal (U10) and meridional winds (V10), and mean sea level pressure (MSLP). Sensitivity experiments explore the use of additional upper-air atmospheric fields; a full list of fields used in this study, their associated symbols used in this manuscript, and the grids on which they are each archived is included in Table 1. The spatiotemporal variations in these fields are considered as well. Spatially, predictors are structured in a forecast-point relative sense. In the control model, GEFS/R forecast values up to \( r = 4 \) grid boxes (\( \sim 2^\circ \)) latitudinally or longitudinally displaced in any direction relative to the forecast point are considered. Temporally, simulated fields are considered at each archive time during the forecast interval, which corresponds to every three hours during the Day 2 period and every six hours during the Day 3 period, for a total of \( t = 9 \) and \( t = 5 \) forecast periods for the Day 2 and 3 periods, respectively. All told, this yields \( t_f (2r + 1)^2 \) model predictors, which yields respectively \( M = 6,561 \) and \( M = 3,645 \) model predictors for the Day Two and Day Three control models. The other category of predictors, background predictors (Table 2), are those which are solely associated with the forecast point, and have no relation to the present meteorology. These include the location of the point, as well as the ARI characteristics of the point and in the surrounding area.
c. Dimensionality Reduction

There are a large number of model predictors, and they are also highly correlated—spatially, temporally, and across variables. With millions of training examples and thousands of features, the forecast problem can become computationally intractable. Further, having many highly correlated features can readily result in model overfitting—making predictions based on noise affecting an individual native feature rather than the underlying signal—a phenomenon commonly termed the “curse of dimensionality” (e.g. Friedman 1997). There are numerous ways these concerns can be addressed; broadly speaking, the most common approaches are either feature selection or feature extraction. In feature selection, a subset of initial predictors are chosen that collectively bear the strongest predictive relationship with the predictand, whereas in feature extraction, a smaller set of new predictors are derived from the original set. Both of these procedures can be performed subjectively through manual means or objectively through automated means. In this case, all of the input predictors are believed to have a physical relationship with extreme precipitation, and choosing only the most predictive fields (e.g. model QPF) and discarding the rest risks removing valuable predictive information not contained in the retained predictor set. The primary issue with the input predictors in this case is not that many may not have any physical bearing on the predictand, but rather that each predictor represents a value at a different point of a continuous field, or a different property at the same point, and are thus necessarily highly correlated to one another. Furthermore, while one could conceivably extract features using field averages or some other pre-determined method, this may not be optimal. For example, it may be better to weight values closer to the forecast point more heavily, while still retaining some information from the far-field predictors. Given the uncertainty in optimally constructing features by manual means, it is more convenient and repeatable to instead extract features objectively. Though it has some
limitations (e.g. Shlens 2014), principal components analysis (PCA; Ross et al. 2008; Pedregosa et al. 2011) is a robust and frequently utilized approach for dimensionality reduction. This creates a small set of uncorrelated predictors that explain the signal in the forecast data and gives insight into the regional modes of atmospheric variability as depicted in the GEFS/R model (explored in more depth in Herman and Schumacher (2018)), while leaving the noise in lower-order principal components (PCs), acting in principle to both alleviate overfitting and manage computational requirements.

d. Machine Learning Algorithms and Sensitivity Experiments

The primary statistical algorithm used in this study is random forests (RFs; Breiman 2001). RFs are in essence an ensemble of decision trees, where traditionally each tree individually makes a deterministic prediction about the outcome of the predictand; the relative frequencies of each possible predictand outcome in the ensemble of trees are then used to make a probabilistic forecast. Much further detail on tree and RF construction and mechanics can be found in Appendix A as well as McGovern et al. (2017) and other sources. There are also several parameters which can be tuned to the particular forecast problem in order to maximize model performance. Four-fold cross-validation is used for model development in this study, whereby each model configuration examined is trained four times, once each on three-quarters of the training data, and then evaluated on the final withheld quarter. To avoid issues of sample independence and approximately mimic information that would be available in an operational context, 974 consecutive initializations are used for each quarter of training data. All parameter settings and sensitivity experiments are evaluated in this framework. The set of RF parameters tuned is described in Appendix A, and the results presented in Appendix B.
In this study, there are a great deal of dynamical model data considered as input information on which the RF can base a prediction. A suite of sensitivity experiments are conducted, as summarized in Table 3, in order to investigate which aspects of forecast information contribute most to forecast skill. Experiments include exploring:

- Sensitivity to the inclusion of horizontal variations in atmospheric fields by varying the previously described predictor radius parameter $R$ from 0 to 4.

- Sensitivity to the inclusion of additional upper-air atmospheric fields by comparing the inclusion and exclusion of two sets of fields as noted explicitly in Table 1. The first incorporates temperature, specific humidity, zonal and meridional winds at 850 and 500 hPa, and 850 hPa vertical velocity in the so-called Upper-Air Core predictor group, while an additional experiment further includes those same fields at 700 and 250 hPa.

- Sensitivity of predictor temporal resolution. Predictor density is three-hourly for Day 2 guidance and six-hourly for Day 3 guidance; models are additionally trained with predictors at twelve-hourly temporal density for both lead times and six-hourly temporal density for the Day 2 forecast model and compared against the control versions.

- Sensitivity to and type the extent of use of ensemble information, a question which has implications for how operational centers allocate their computational resources. Using forecast information from only the GEFS/R’s control member in model training (CTRL) is compared with using the ensemble median from the full ensemble (MEDIAN), and then further with the use of the ensemble second-lowest and second-highest values for each atmospheric field in conjunction with the median (CNFDB) to evaluate the impact of this dimension of forecast information, following the findings of Herman and Schumacher (2016b), which found relatively little sensitivity in performance with respect to how ensemble information is used, but
using the near-minimum, median, and near-maximum values outperformed using the mean
and spread.

- Sensitivity to predictor pre-processing methodology. Models are trained with and without
the aforementioned PCA pre-processing step, and an assessment of the effect of this pre-
processing step on model skill is made by comparing the two.

- The effect of region size on forecast skill, hypothesizing models trained for larger regions may
exhibit higher skill due to more available training data. This is performed by aggregating the
ROCK and SW regions into a new WEST one, combining the Southern Great Plains (SGP),
Northern Great Plains (NGP), and Midwest (MDWST) regions into a CENTRAL region,
and collecting SE and NE regions into a single EAST region, while leaving PCST—with its
unique extreme precipitation climatology—unperturbed.

- Sensitivity of model performance as a function of model algorithm, specifically by comparing
with logistic regression (LR), a common and comparatively simpler alternative to statistically
deriving forecast probabilities. Further discussion of LR and other machine learning alterna-
tives to the RF algorithm is included in Appendix A.

**e. Model Evaluation**

Based on the parameter tuning and sensitivity experiment results, final model configurations are
selected. The final model is run over a completely withheld 4-year evaluation period spanning
September 2013–August 2017. The forecasts generated from the final model are compared with
those from the full ensemble of raw GEFS/R QPFs, as well as the full 50-member ECMWF global
ensemble, accessed from TIGGE (Molteni et al. 1996; Bougeault et al. 2010). The comparison
with the former provides an assessment of what improvement, if any, these models yield compared
with the raw guidance from which their forecasts are derived when evaluated in a real-time setting.

The latter, meanwhile, provides an assessment for how these forecasts compare with state-of-the-
science operational ensemble guidance available at these lead times. To make these comparisons,
the QPF from each ensemble member of the two ensembles is regridded onto the ∼4.75km Stage
IV HRAP grid on which the Atlas thresholds lie using a first-order conservative scheme (Ramshaw
1985). These regridded QPFs are then compared with the 1-year and 10-year ARI thresholds to
create deterministic exceedance forecasts with respect to the two thresholds for each ensemble
member. These binary grids are then upscaled to the GEFS/R grid using the same procedure as
the verification upscaling: any exceedance in the downscaled grid corresponds to an exceedance
at the nearest GEFS/R point in the upscaled grid. Since the predictand categories are necessarily
mutually exclusive, the 1-year ARI exceedance grids are modified so that any member forecasting
a 10-year ARI exceedance at a point is not forecasting a between 1-and-10-year exceedance at that
same point and time period. The prevailing operational method of generating forecast probabili-
ties from a dynamical ensemble—democratic voting, whereby the fraction of ensemble members
forecasting the event is used as the forecast probability (e.g. Buizza et al. 1999; Eckel 2003)—is
applied to each ensemble to generate the exceedance probabilities for the reference forecasts.

Skill, both in the final assessment of model performance as well as in all aforementioned sen-
sitivity experiments, is quantified by means of the Rank Probability Skill Score (RPSS) with a
climatological reference:

\[
RPSS = 1.0 - \frac{\sum_{d=1}^{D} (\sum_{p=1}^{P} (\sum_{m=1}^{K} (\sum_{j=1}^{P} P_{jpd} - O_{jpd})^2)))}{\sum_{d=1}^{D} (\sum_{p=1}^{P} (\sum_{m=1}^{K} (\sum_{j=1}^{P} P_{clim_{j}} - O_{jpd})^2)))}
\] (1)

with D forecast days; P forecast points; K predictand categories; \(P_{jpd}\) and \(O_{jpd}\) corresponding
respectively to the forecast probability and observance of predictand category j on day d and at
point p; and \(P_{clim}\) corresponding to the climatological frequency of occurrence, as defined by
the respective ARIs of the predictand. A score of 1.0 indicates a perfect forecast, and a score of 0.0 indicates model performance equivalent to forecasting climatology. Final assessment also includes analysis of reliability, both subjectively through reliability diagrams, and quantitatively via the Murphy (1973) decomposition of the Brier score (BS) for category $j^*$:

$$BS_{j^*} = \sum_{n=1}^{N} (P_{N,j^*} - O_{N,j^*})^2 = \frac{1}{N} \sum_{c=1}^{C} N_{c,j^*} (P_{c,j^*} - \overline{O_{c,j^*}})^2 - \frac{1}{N} \sum_{c=1}^{C} N_{c,j^*} (\overline{O_{j^*}} - \overline{O_{c,j^*}})^2 + \overline{O_{j^*}} (1 - \overline{O_{j^*}})$$

(2)

where there are $N = DP$ total forecasts, broken into $C$ discrete probability bins with $N_c$ forecasts being issued for each bin $c$. $\overline{O_{j^*}}$ denotes the climatological (based on the period of record) frequency of observing event category $j^*$ and $\overline{O_{c,j^*}}$ denotes the proportion of forecasts in probability bin $c$ observing event category $j^*$, where $j^*$ is the aggregation of event categories of at least $j$ in the RPSS framework. $\overline{O_{j^*}} (1 - \overline{O_{j^*}})$, the so-called “uncertainty” term, also represents the BS of a climatological forecast. Converting to a Brier skill score (BSS) framework by dividing out by this term:

$$BSS_{j^*} = 1.0 - \frac{BS_{j^*}}{BS_{clim,j^*}} = \frac{1}{N} \sum_{c=1}^{C} N_{c,j^*} \left( \frac{(\overline{O_{j^*}} - \overline{O_{c,j^*}})^2}{\overline{O_{j^*}} (1 - \overline{O_{j^*}})} \right) - \frac{1}{N} \sum_{c=1}^{C} N_{c,j^*} \left( \frac{(P_{c,j^*} - \overline{O_{c,j^*}})^2}{\overline{O_{j^*}} (1 - \overline{O_{j^*}})} \right)$$

(3)

This analysis is conducted for both the 1- and 10-year thresholds.

Skill calculations and comparisons are made for the host of sensitivity experiments and for each region, lead time, and model configuration. For each comparison, statistical significance is assessed by bootstrapping to obtain identical sets of cases for each of the two forecast sets being compared. Skill scores are derived from the subsample of each forecast set, and a skill difference is computed. This process is repeated 1000 times to generate a distribution of skill differences, and statistical significance is ascertained with respect to whether the 0.5th and 99.5th percentile skill score difference values from the bootstrap trials overlap zero. This 99% confidence bound is used in contrast to 90% or 95% bounds to compensate for concerns arising from conducting stas-
tical significance analysis on numerous different comparisons. While some uncertainty analysis is included in the figures presented, much of the statistical significance difference results discussed in-text are omitted for the sake of concision.

3. Results: Sensitivity Experiments

Examining forecast skill as a function of time step between atmospheric field predictors (i.e. the CORE_LTIME models of Table 3; Fig. 4a), two striking findings concern 1) the large variations in forecast skill across regions and 2) the evidently low sensitivity of forecast skill to time step length within any given region. For the 3-hour time step, predictors are gathered from a total of 9 forecast times; with the 6-hour step, 5 forecast times are used; and with the 12-hour time step, a total of 3 forecast times are used. The 12-hour time step therefore has one-third the total number of predictors as the model with the 3-hour time step, but still yields nearly identical forecast skill results. In most regions and forecast periods, there is a slight degradation in performance going from the 6- to 12-hour time step, but the difference is not generally statistically significant by a 99% bootstrap skill score difference test (not shown). The one exception to this is in the PCST region, which has much higher skill overall than the other regions for both forecast periods, and exhibits somewhat higher sensitivity to the predictor time step than the other regions, particularly in going from 6-hours to 12-hours, with RPSS differences of approximately 6%.

Similar to the temporal resolution findings, there is a general lack of sensitivity as a function of predictor spatial extent (Fig. 4b). This finding comes in stark contrast to that of Herman and Schumacher (2016b), which found great sensitivity of predictor spatial extent in forecasting airport flight rule conditions. Albeit weak, a slight improvement in skill for most forecast period, region combinations can be noted with increasing predictor radius, often to the extent that the skill difference between 0 and 4 grid box radii is statistically significant (not shown). Two regions in
particular, the NE and PCST, exhibit by far the most sensitivity to predictor spatial extent, with differences of roughly 0.02 observed over the evaluated interval. Also of note is that a radius of 4 grid boxes—the highest number evaluated—did not always yield the best performance results; most notably, the Day 2 model for the NE region maximized skill at a radius of 2, with a slight deterioration of forecast skill with increasing radius thereafter. In those regions where the GEFS/R cannot explicitly resolve the processes responsible for producing extreme precipitation, the RF is ultimately making forecasts more on environmental factors; these do not vary drastically in time or space, and thus a single number or small set of numbers at or immediately surrounding the forecast point are sufficient to characterize the basic properties of the environment. This is all that the RF is really using for much of its predictions (see Herman and Schumacher 2018, for more detail). However, in regions impacted more readily by larger scale systems where the dynamical model can more directly simulate the precipitation processes such as PCST and the NE, the spatial variations in atmospheric fields carry more signal rather than noise and thus contribute more predictive value.

Like varying spatial and temporal density, there is relatively little sensitivity to the inclusion of more atmospheric fields (Fig. 5a). Slight but consistent improvement is observed in adding the core upper-air fields as predictors, but adding further levels beyond the core group was found to not improve predictive skill, and actually resulted in a decrease in skill for the PCST, NE, ROCK, and SE regions—those which are most affected by larger scale precipitation systems. Though still rather small, somewhat more distinct sensitivity to type of ensemble information included (Fig. 5b) can be seen here across all regions, with improvements seen using predictor information from the GEFS/R ensemble median versus using only the control member, and slight further improvement using the ensemble second-from-minimum and second-from-maximum in addition to the ensemble median. The largest differences in magnitude are again for the PCST.
region, but in this experiment, clear and statistically significant (not shown) improvements are also seen for low skill, convectively active regions such as MDWST.

Aggregating regions (Fig. 5c) results in a slight degradation in forecast skill. In principle, it is possible for a decision tree to automatically forecast for specific regions by splitting first on the latitude and longitude predictors, and then further partitioning based on meteorological variables thereafter. However, these findings demonstrate that there is some—albeit limited—utility in manually partitioning training data with distinct hydrometeorological relationships, rather than relying on the machine learning algorithm to discern the distinction automatically. Comparing the impact of applying PCA pre-processing to the RF (Fig. 5d, leftmost two columns), performing PCA tends to either improve performance, as is the case for the PCST, NE, SW, and MDWST regions, or make little difference, as seen in the ROCK, NGP, SGP, and SE regions. The positive differences tend to be larger in magnitude, both in relative and absolute senses, for Day 2 model versions compared with Day 3. Forecasts produced through LR tend to be substantially worse than those generated by RFs (Fig. 5d, center columns). However, the exact magnitude to which this is the case varies by region; substantial differences in skill are seen between RF and LR forecasts for the SW, ROCK, and SGP regions, while there is almost no skill difference between the Day 3 forecasts in the PCST region. This may suggest the linear assumptions inherent to the LR algorithm perform better in larger scale systems than in the more convectively active ones in which the responsible processes are highly nonlinear, but this causality is not entirely clear. Finally, a weighted average of RF and LR forecasts outperforms its component members for all regions and forecast periods. The extent of overperformance is strongly tied to the skill difference between the RF and LR models; when the skill difference is small, the value of the weighted average is comparatively large to when the RF performs much better than LR (cf. Fig. 5d PCST and SW lines). Since these weighted averages performed the best in cross-validation, a weighted
average using each of the CTL_NPCA, CTL_PCA, and CTL_LR models was chosen for the final model configuration.

4. Results: Final Model Performance

For both the final ML models and the forecasts from the raw QPFs of both the GEFS/R and ECMWF (Fig. 6), a usually statistically significant deterioration in forecast skill from Day 2 to Day 3 is evident in each CONUS region over the four year test period. Forecast skill is significantly higher in regions with extreme precipitation associated partially or primarily with synoptic scale precipitation episodes, such as PCST, SW, and ROCK, rather than smaller-scale convective systems that characterize extreme precipitation as in the NGP, SGP, and MDWST regions. At an extreme, the NGP and SGP GEFS/R raw QPFs exhibit no skill in predicting ARI exceedances at these lead times. Especially for the ML models, the bigger Day 2 vs. Day 3 skill differences are also seen where the skill is higher, again suggesting the direct forecasting of the precipitation as opposed to forecasts more reflecting the forecast environment, either dynamically via parameterized convection in the case of raw QPFs, or directly in the case of the ML model forecasts. Furthermore, the ML models exhibit a larger skill deterioration between Days 2 and 3 than either of the raw ensemble forecast sets.

Comparing the forecast systems, the ECMWF forecasts consistently and statistically significantly outperform the GEFS/R forecasts at all lead times except in the SE region (Fig. 6). Encouragingly, the ML model forecasts are statistically significantly more skillful for all eight regions and both lead times compared with the GEFS/R forecasts from which they are based. The post-processing is thus clearly accomplishing its purpose of improving forecast skill. But it is also apparent that the GEFS/R is not a state of the science model for extreme QPF prediction given its lower skill compared with the ECMWF. The real test of the ML model then is how it compares...
with current best operational guidance for these lead times, represented here with the ECMWF ensemble. The comparison (Fig. 6) is generally quite favorable, with the Day 3 ML forecasts outperforming even the Day 2 ECMWF forecasts across all regions except ROCK and PCST. In the non-western regions, the extent of overperformance is quite considerable when comparing equal lead times, with skill score improvements of factors of two to three seen in many comparisons. In the ROCK and PCST regions, the ML and ECMWF forecasts performed about equally at Day 2, and ECMWF performed slightly better at Day 3. Overall, the ML models demonstrated ability to consistently outperform current operational model guidance, especially in convectively active regions where there is no operational guidance that can dynamically resolve the physical processes producing extreme precipitation at these lead times.

Reliability diagrams of Day 2 raw GEFS/R and ECMWF forecasts (Fig. 7) reveal highly overconfident probabilistic exceedance forecasts for all regions, both severity levels, and both ensembles as evidenced by the shallow slope relative to the one-to-one line in each panel. The raw GEFS/R forecasts (Fig. 7a,b) are relatively sharp, with more than 0.01% of forecasts falling into each probability bin above 10%, and a vast majority of zero probability forecasts (not shown). For all regions, there are cases where every ensemble member has simultaneously predicted a 1-year exceedance (Fig. 7a), but the same is not true for 10-year exceedance predictions in the northeastern regions: NE, NGP, and MDWST (Fig. 7b). The ECMWF (Fig. 7b,d) is also overconfident, but we see that it is also negatively biased for all cases. Its degree of overconfidence is dampened compared with the GEFS/R, and it is not as sharp, with fewer very occurrences of very high forecast probabilities except in the westernmost regions of ROCK and PCST (Fig. 7b, inset panels). With 50 members rather than 11, there is also substantially more resolution across the probability spectrum in the ECMWF forecasts. By the very nature of how these forecasts are generated, quite a bit of sharpness is inherent at the cost of reliability, since it is not possible for probabilities near
the climatological event frequency to be issued for either raw ensemble, but particularly for the
GEFS/R.

The Day 2 reliability diagrams for 1-year exceedance forecasts from the different components of
the final model—CTL_{NPCA}, CTL_{PCA}, and CTL_{LR}—are shown in Figure 8. The CTL_{NPCA}
(Fig. 8a) shows markedly different characteristics than either of the raw ensembles. In particu-
lar, all of the regions exhibit an underconfidence signal, with low probability events below about
2% for 1-year events (Fig. 8a) occurring with observed relative frequencies below the forecast
probabilities. The relative event frequencies are conversely appreciably higher than the forecast
probabilities would indicate for probabilities above 5%. Among the regions, the PCST probabili-
ties are the most negatively biased, while NE probabilities are the most positively biased. Overall,
reliability is much better than for either raw ensemble, but this comes at the expense of sharpness.
Less than 1 in 10,000 forecasts are above about 20% for (Fig. 8a, inset panels), and maximum
probabilities are in the 30–80% range depending on the lead time and region, compared with
100% for all lead times and regions in the raw ensembles. The CTL_{PCA} model (Fig. 8b) exhibits
very similar reliability characteristics to the CTL_{NPCA} model, including the underconfidence,
reduced sharpness compared with the raw ensembles, and different regional probability bias char-
acteristics. It tends to be more negatively biased than CTL_{NPCA} at low and high probabilities (cf.
Fig. 8a,b), correctly so at high probabilities and undesirably so at low ones. The CTL_{LR} model
(Fig. 8c) exhibits some similarities and some differences with the RF-based models. PCST fore-
casts are consistently the most negatively biased, followed by ROCK and the SE, with NE region
forecasts being the least negatively biased. However, unlike the RF-based forecasts, the LR model
issues more high probabilities; for example, forecasts in the highest probability bin were issued for
most regions (Fig. 8c). At the highest probabilities, the forecasts revert to being positively biased,
as they are for events with probabilities issued in the 0.01–1% range. At very low probabilities ,
LR-based forecasts are substantially more negatively biased than for RF-based forecasts, leading to considerable overconfidence overall when considering that the vast majority of forecasts issued occur on this low probability end of the spectrum. While LR (and regression in general) is effective at removing bias in a global sense, since a single regression equation must necessarily apply globally to all forecasts, it inherently cannot perform more localized, context-depend forms of bias correction, leading to forecast probability-dependent model biases.

The final ML model reliability (Fig. 9) unsurprisingly reflects a blend of the component members, retaining some of the underconfidence of the RF-based models while adding a bit of sharpness from the CTL_LR model in regions where it verified skillfully enough in cross-validation (e.g. PCST, Fig. 5d) to garner much weight. The probability distribution for 1-year exceedance events is not markedly different between the Day 2 and Day 3 forecasts (cf. Fig. 9a,c), but the relatively higher probabilities issued for 10-year exceedances in Day 2 do not occur at the Day 3 lead times (cf. Fig. 9b,d). This is consistent with increasing confidence in very extreme events with decreasing lead time—something seen very pronounced in the final model, but to a much lesser extent in the raw ensemble forecasts.

The relationship between the reliability analysis and skill via the Brier score decomposition (Murphy 1973) quantitatively solidifies many of the general observations discerned by inspection of the reliability diagrams. Though sharper than competing forecasts, the raw GEFS/R forecasts consistently exhibit the worst resolution component contribution to forecast skill for all regions and severity levels, both for Day 2 forecasts (Fig. 10a,c) and Day 3 forecasts (Fig. 11a,c) due to an inability to actually distinguish events from non-events by resolving the responsible physical mechanisms. The final ML models exhibit better resolution term skill contributions than the ECMWF ensemble forecasts, with the exception of the ROCK and NGP regions for 1-year events (Fig. 10a, 11a). Between the component models, resolution term skill tended to best for
CTL-NPCA forecasts over the test period, particularly at the 10-year severity level (e.g. Fig. 10c) but the extent of the difference tended to be relatively small and there were numerous instances where PCA-based models exhibited more resolution. The weighted average consistently exhibited higher resolution than any of the component members. With respect to the reliability contribution to skill (Fig. 10b,d Day 2; Fig. 11b,d Day 3), ECMWF forecasts were, perhaps surprisingly given the lack of explicit calibration, the most reliable forecast set for all regions and lead times, while in many cases the ML models had a more negative contribution to the total skill than the raw GEFS/R, likely resulting from the underconfidence. The resolution term is at largest one and at least zero in this decomposition, while the reliability term is at most zero. The magnitude of the resolution terms is consistently several factors larger than the reliability term for all forecast sets, and the differences in that term generally have a larger absolute impact on the overall Brier skill scores.

Lastly, while by no means a comprehensive characterization of the system, a sample of real cases over the test period are presented to illustrate some of the strengths and weaknesses of the system. On the evening of 19 May 2015 and morning of 20 May 2015, a vigorous mesoscale convective system developed over southern Oklahoma and northern Texas, producing very heavy rainfall that contributed to historic flooding in the region during May 2015 (e.g. Wolter et al. 2016). Stage IV analysis (Fig. 12a) reveals that the 24-hour precipitation totals exceeded 1-year ARI thresholds within much of an E/W band encompassing the region, with embedded areas of 10-year exceedances along the state border region (Fig. 12b). While the ECMWF ensemble forecasts indicate some possibility of extreme precipitation in that region during this time frame at Day 3 (Fig. 12d), the probabilities are displaced too far to the south and west, and the probabilities of 10-year exceedances are very low. There is some improvement in positioning with the Day 2 forecast (Fig. 12c), but it remains too far west and with probabilities still quite low, particularly
at the 10-year ARI level. Raw GEFS/R forecasts at Day 3 (Fig. 12f) indicate quite high risk for a 1-year exceedance over a fairly narrow area, better positioned than the ECMWF ensemble at the same lead time but still too far to the west. Outside of this area, the GEFS/R indicates almost no risk of an extreme rainfall event, and also indicates no risk of a 10-year exceedance anywhere in the domain. The Day 2 forecast (Fig. 12e) looks similar to the Day 3 outlook, except that the probabilities are reduced somewhat in the target area, which also has incorrectly displaced further to the south and west. The ML model depicts a much different picture. It exudes much less confidence, with lower maximum probabilities compared with either raw ensemble, but non-zero exceedance probabilities of both 1- and 10-year exceedances across much of the domain for both Days 3 (Fig. 12h) and Day 2 (Fig. 12g). Importantly, the model elevated probabilities compared with the raw guidance in the place that extreme precipitation was actually observed (to the east of where it was forecast in the GEFS/R). In fact, at Day 2 (Fig. 12g), the probability maximum is located right where the heaviest precipitation actually occurred, displaced well to the north and east of where it was forecast in the GEFS/R (Fig. 12e). Additionally, while still low, the 10-year event probabilities are much higher over the verifying area when compared with either raw ensemble, with maximum Day 2 probabilities of around 30% and 3% for 1-year and 10-year exceedances, respectively. Finally, in contrast to the raw guidance, the ML model became increasingly confident in an event occurring with decreasing lead time (cf. Fig. 12g,h).

A different mesoscale precipitation produced extreme precipitation over southwestern Wisconsin, southeastern Minnesota, and northeastern Iowa during the evening and overnight hours of 21 September and 22 September 2016, respectively. Based on ST4 QPE (Fig. 13a), much of the area experienced 1-year ARI exceedances for the 24-hour period ending 1200 UTC 22 September 2016, and within the 1-year exceedance area, there were many embedded cells that produced 10-year ARI exceedances (Fig. 13b). ECMWF forecasts at Day 3 indicated risk of extreme rain-
fall, even at the 10-year severity level (Fig. 13d), but the location was poor, with exceedance probabilities high in eastern Minnesota and northern Wisconsin where extreme rainfall was not observed, and very low probabilities in northeastern Iowa and southeastern Wisconsin where it was. Both the positioning and risk of very extreme precipitation improved for the Day 2 forecast issuance (Fig. 13c), but probabilities still remained too far to the north. The GEFS/R at Day 3 (Fig. 13f) indicated very little risk of extreme precipitation in the area, with just one member correctly predicting a 1-year exceedance in southeastern Minnesota. The risk of an event occurring within the domain increased for the Day 2 issuance, but the locations got worse, with maximum risk indicated in eastern Nebraska, western Iowa, and northeastern Wisconsin, with the only 10-year prediction occurring in the latter location. Somewhat like the raw GEFS/R, the ML model had only some indication of extreme precipitation risk at Day 3 (Fig. 13h). However, it both had the higher probabilities (near 10% in both cases) distributed over a much larger area, and indicated some risk of a 10-year event, with probability maxima near 1.5%. Additionally, it had the maximum probability axis nearly collocated with where heaviest precipitation occurred, well to the south of the ECMWF probabilities, albeit still slightly too far to the north. The Day 2 forecast issuance (Fig. 13g) was largely similar. The two main changes are a correctly increased risk in the area where the event actually verified, and an incorrectly increased risk of heavy precipitation in eastern Nebraska where the raw GEFS/R had heavy precipitation on Day 2 (Fig. 13e).

5. Discussion and Conclusions

An ML model based on RFs and LR is used to generate CONUS-wide probabilistic forecasts for the exceedance of 1- and 10-year ARI thresholds for 24-hour precipitation accumulations during the Day 2 and Day 3 periods. Approximately eleven years of GEFS/R forecasts, in particular the ensemble median, are used to train these models, and forecasts are made using numerous
simulated atmospheric fields (Table 1) varying in both space and time, in addition to a variety of geographic and climatological forecast predictors (Table 2). Separate models are trained for each of the two 24-hour periods and for each of eight different regions of CONUS, as depicted in Figure 3. A variety of sensitivity experiments are performed, as outlined in Table 3, to ascertain the utility of different aspects of forecast information in predicting locally extreme precipitation. Finally, the final forecast models were evaluated, and compared with forecasts based only on the ensemble of raw QPFs from the GEFS/R and ECMWF. The ML models trained in this study demonstrably outperformed the raw GEFS/R forecasts for all regions and forecast lead times (Fig. 6), often more than doubling the forecast skill and adding substantially more than 24-hours lead time improvement in forecast skill. With the exception of the PCST and ROCK regions, the same held for comparison of the ML model forecasts with ECMWF ensemble forecasts as well. Both raw ensembles tended to be negatively biased and highly overconfident in predicting extreme QPFs (Fig. 7), particularly at the 10-year ARI for central CONUS regions; this was reversed in the final ML model forecasts, which were more reliable at higher probabilities, but generally underconfident (Fig. 9).

In general, unlike past studies (e.g. Herman and Schumacher 2016b), in most regions, the temporal resolution and extent of spatially displaced predictors from the forecast point considered had little to no impact on forecast skill (Fig. 4), in addition to the use of upper-level information and additional ensemble information (Fig. 5). These results are suggestive of two findings. First, most of the relevant information about predictors displaced spatiotemporally from the forecast point, other atmospheric fields, or other ensemble member information, can be derived with at least moderate accuracy using just the information from the ensemble median from a group of core set of fields collocated and concurrent with the forecast; that is, these additional predictors contain only limited independent forecast information, at least for this coarse dynamical model and this under-
dispersive ensemble configuration. It also suggests that, for the most part, the predictive ability is coming primarily through a characterization of the overall environment, which can be reasonably summarized with only a subset of predictors, rather than the simulated spatiotemporal variability and full 3-D characterization of the atmospheric evolution in the underlying dynamical model. This finding comes in contrast to similar studies of other forecast problems using the GEFS/R, such as the Herman and Schumacher (2016b) study which investigated using the GEFS/R to create ML-based probabilistic forecasts of cloud ceiling and visibility at different airports and found considerable value in the inclusion of spatially displaced predictors. However, there is at least one major exception; none of this really held for the PCST region; here, more complex models with more predictors did notably improve forecast skill. This is perhaps in part because the physical processes associated with extreme precipitation are much better resolved in the GEFS/R in this region compared with the others, and so the added information adds usable forecast utility beyond simply duplicatively characterizing the atmospheric environment for the forecast. The largest skill difference of the sensitivity experiments came for most regions in changing algorithmic assumptions and processes (Fig. 5d); the simpler linear assumptions of LR tended to degrade forecast skill compared with the more limited assumptions underlying the RF models.

The results of this study reveal that the application of more sophisticated statistical methods and ML algorithms such as RFs can demonstrably improve forecasts of extreme precipitation and potentially other rare, high-impact weather events in the medium range when compared with the methods and techniques that are most prevalent in forecast operations today. One unique aspect here is the scope of this model; while most past studies which employed these techniques for numerical weather prediction have focused on a small domain, or just a sampling of points, the models trained here demonstrate an ability to generate skillful, reliable forecasts year-round for all of CONUS and a range of lead times. There are many forecast problems that remain to be
explored, but the results of this study and others strongly suggest that further development and application of these data-intensive statistical techniques could substantially improve our forecasts over the current state of the art, even compared with using more sophisticated dynamical models. To that end, implementation of this methodology for operational use to assist Weather Prediction Center forecasters with the development of their excessive rainfall outlooks is currently underway.

This forecast technique presents some advantages over purely dynamical approaches, as dynamical models are inherently limited by two factors by which these statistical techniques are not. First, dynamical models require ever increasing computational resources for increasing model resolution; constraints on computing power prevent sufficient resolution to directly resolve many small-scale processes, many of which are observed in the highest impact weather phenomena. Second, dynamical models are limited by our physical understanding of the processes we are attempting to simulate or forecast. Machine learning algorithms, in contrast, can detect predictive patterns in the available information even in places where we do not know or understand the physical connection between the information and the phenomenon which we wish to predict. While they are also limited in complexity by computational and data resources, the strict limits on resolvability are not there: physical resolution can often be gained through post-processing of larger scale information. There is thus ample reason to believe that further investigation of these techniques for NWP is a worthwhile venture, and eventual implementation into forecast operations could help forecasters with their tasks by skillfully synthesizing many different sources of forecast information to help alleviate their often time-pressed schedules. This in turn can aid end-user preparedness and, in the case of high-impact events, hopefully help to protect lives and property.

One of the main advantages of the methods explored in this study compared with other popular machine learning methods, in addition to their computational tractability, is the ability to visualize their output and gain insights into detecting and quantifying specific biases in the underlying
GEFS/R model, and physical insights into the most valuable forecast information for predicting locally extreme precipitation. For reasons of focus and brevity, the diagnostics that shed light on these insights have been omitted from this manuscript and are presented instead in a companion paper focused on the diagnostics (Herman and Schumacher 2018) rather than the forecasts and forecast process explored in depth here.

Some limitations of this work are worthy of note. Stage IV precipitation is used as truth for this study; though there is not a clearly better verification source available, it does have its drawbacks. It does have some spurious quality control issues, and often struggles in areas of complex terrain due to radar beam blockage, interference, and limited gauge coverage (Herman and Schumacher 2016a; Nelson et al. 2016). Since the model is trained to forecast Stage IV QPE exceedances, this can lead to some idiosyncrasies and other anomalies associated with the biases observed in the Stage IV product. One such anomaly is the persistent presence of very small areas of exceedances in some regions of complex terrain during times of favorable convective conditions. This can be removed by quality control procedures to some extent, but some artifacts do remain. This happens most prominently in the terrain of western New Mexico; a small region there has many more instances of ARI exceedances over both the training and test periods than any other part of CONUS. The ML-based models recognize this, and for the SW region consistently issue much higher probabilities in this region. In one sense, this is correct—it is correctly predicting what it was trained to predict—but is still undesirable behavior due to a disparity between “truth” in the study and the true extreme rainfall risk. Solutions to this issue and related issues in other parts of the country must be explored in order to maximize operational utility. Additionally, while the choice of using the ARI framework was an intentional decision and provides numerous benefits, it is not an end-all for predicting heavy precipitation impacts. While ARIs often have better correspondence with impacts than a fixed threshold, there are still regional discrepancies in which ARIs
have optimal association with impacts, and the framework employed here does not account for antecedent conditions, which can be critical for assessing flash flood risk. More investigation into the relationship between QPE exceedances and rainfall impacts should be performed to maximize the practical significance of the model predictand.

Additionally, the predictors for this study come from a very coarse and otherwise rather antiquated global model. The GEFS/R was used for this study because, unlike almost any other dynamical model, it has been nearly static for a very long period of record and has nearly stationary bias characteristics—an essential property for performing this kind of analysis. However, the models trained herein are not working off of the ‘state of the art’ of flash flood predictors. The longer range Day 2 and Day 3 lead times were chosen for this study in part because the discrepancy between GEFS/R forecast quality and ‘state of the art’ is smaller at these longer lead times due to less convection-allowing guidance being available, and higher-resolution models degrading in utility with increasing forecast lead time (e.g. Zhang et al. 2003, 2007).

There are also some complications that must be considered for real-time implementation. As one example, the regional models are trained completely independently of one another, with different training data and different solutions. Consequently, they can occasionally give rather different predictions on nearly identical inputs, resulting in undesirable probability discontinuities across region boundaries. Appropriate methods for removing probability discontinuities in space must be further explored.

Future work will seek to alleviate these limitations in a variety of ways. Exploration of using different predictands, likely combining hydrometeorological information from a variety of sources, will be made for more explicit flash flood prediction. This may involve a regionally varying predictand definition, with some ARI thresholds better corresponding to flash flood impacts in some regions compared with others. Additionally, although a large number of predictors were explored...
in this study, there are many additional choices for predictors that could ostensibly further improve forecast skill. While atmospheric fields are represented here in absolute terms, it may be beneficial to instead represent some fields relative to the local climatology of the forecast point in terms of standardized anomalies. This is particularly true for fields like PWAT, where standardized anomalies have often shown better correspondence with precipitation impacts across varied regions than absolute values (e.g. Junker et al. 2009; Graham and Grumm 2010; Nielsen et al. 2015). More exploration of derived fields of physical relevance to extreme precipitation processes should also be explored. Some possible examples include upslope flow to gauge forcing for ascent by the horizontal wind, column mean wind to ascertain potential for slow-moving storms, and deep-layer shear as a metric for supercell potential.

This study also focused on a rather specific time interval and took all dynamical predictors from a single, somewhat antiquated ensemble system. Future expansion both to the 12–36 hour Day 1 period and beyond the Day 3 period will be explored, including predictors from more contemporary CAM guidance and potentially including observations as well for the shorter lead time forecasts. Operational models also tend to undergo periodic upgrades and thus do not remain static like the ensemble system used here. The sensitivity of ML model performance to changes in dynamical model bias characteristics that result from these upgrades is a question of considerable operational relevance and an additional factor worthy of future investigation. It was also seen that the ML models suffered to varying degrees from underconfidence and, in some instances, negative bias. Methods of probability calibration of the ML model probabilities as a final post-processing step (e.g. Hagedorn et al. 2008; Hamill et al. 2008; Bentzien and Friederichs 2012; Herman and Schumacher 2016b) should be explored in future work, and parameter choices reconsidered in light of this additional calibration. Finally, this study only explored a subset of available machine learning algorithms. Other choices, including adaptive learning algorithms, may be able to better
exploit predictor-predictand relationships, appropriately update to reflect changes in an underlying
dynamical model, and produce superior forecasts for the locally extreme precipitation and flash
flood forecast problem (e.g. Liu et al. 2001; Roebber 2015; Pelosi et al. 2017).

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APPENDIX A

Algorithm Descriptions

a. Random Forests

As noted in the main text, RFs are simply an ensemble of decision trees. Decision trees consist
of a network of two types of nodes: decision nodes and leaf nodes. Decision nodes each have
exactly two children, which may be either decision nodes or leaf nodes, with a binary split based
on the numeric value of a single input predictor determining whether to traverse to the left or right
child. A leaf node has no children and instead makes a categorical prediction of the outcome of
the input example based on the leaf’s relationship to its ancestor nodes. For a given forecast, one
begins at a decision tree’s root, traversing through its children based on the relative value of the
forecast’s predictors to each decision node’s threshold critical value for the predictor associated
with the node. This process is repeated until a leaf node is reached; its value corresponding to the
leaf becomes the tree’s deterministic prediction.

Decision trees can be a powerful approach for a wide array of applications, but they also have
several significant drawbacks. In particular, they are very prone to overfitting (e.g. Brodley and
Utgoff 1995), fitting to the noise of the training data rather than just the underlying relationships.
They also don’t convey any information about forecast uncertainty, as would be the case in a prob-
abilistic framework. RFs are used instead to alleviate these concerns by producing a probabilistic
forecast in a way that can significantly decrease error from overfitting the supplied training error
with only a slight increase to error from oversimplistic model assumptions, provided the trees are
sufficiently uncorrelated. The difficulty then revolves around generating a large set (forest) of
skillful decision trees that are not strongly correlated. The decision tree generating procedure de-
scribed above is deterministic: a given set of training data will always produce the same decision
tree. A forest of identical decision trees, of course, adds no value over using a single decision tree.
Two additional processes—tree bagging and feature bagging—are employed to produce unique
trees. Tree bagging produces unique trees through a straightforward bootstrapping procedure.
Specifically, a forest of size B is formed from the n training examples by creating B samples of
size n, with replacement, from the original training data, and running the decision tree algorithm
on each sample. Overfitting due to correlated trees can still occur under this approach, particularly
if a small subset of the original features are much more robust predictors of the verifying category
than the rest (Breiman 2001; Murphy 2012). To overcome this problem, feature bagging is also
employed, whereby only a random subset of the m original input predictors are considered at each
decision node; the size of the random subset is denoted here as S; 1 ≤ S ≤ m. This combination can
result in a set of B largely uncorrelated trees, each of which is individually fairly skillful.
With any machine learning algorithm, there are numerous considerations in the actual model construction, which manifest themselves in tunable parameters. Compared with other machine learning algorithms, such as gradient boosting or support vector machines, RFs are often praised for their relative insensitivity to their parameters with respect to model performance, but it is nevertheless important to explore the parameter space in order to realize the full utility of the algorithm. The forest size $B$ is perhaps the most obvious parameter. The general relationship between model performance and $B$ is well known and consistent across all prediction problems; it starts quite low at very low $B$, initially increases rapidly with increasing $B$, and then slowly asymptotes to some threshold performance limit as the relationships between input features have been fully explored by the forest and the inclusion of new trees becomes redundant. Larger forest sizes require more computational expense, so the goal is to select $B$ such that it is small enough to be computationally tractable but large enough to be near the performance limit. Another parameter noted above is $S$, the number of features to consider at each node split. If this number is too small, model performance may suffer from only considering irrelevant or otherwise unpredictive features in the context of the node; if $S$ is too large, performance will also suffer because of underdispersive trees producing an overfit forest solution. Another frequently explored parameter is the splitting criterion evaluation function. Most commonly used are either the Gini impurity or the information gain; past studies have shown that this choice is not important for many forecast problems. Information gain is used in this study; it can be expressed for a training set $T$, candidate splitting feature $x_a$ and candidate split value $v_a$ as:

$$IG(T, x_a, v_a) = H(T) - H(T | x_a < v_a)$$  (A1)
where \( H(T) \) is the so-called entropy of a tree, defined for each of the \( K \) verifying categories, with each category \( i \) having forecast probability \( p_i \), as:

\[
H(T) = - \sum_{i=1}^{K} p_i \log_2 p_i
\]  

(A2)

The chosen splitting feature and split value are selected among those considered which maximize Equation A1 (e.g. Quinlan 1986; Murphy 2012). However, there are two other parameters that have the most substantial influence on model performance. The first, denoted \( Z \), is the minimum number of training examples required to split a node. Traditionally, RFs create a leaf only once a node is ‘pure’, that is, all the remaining training examples associated with that node have the same labels (event outcomes). In this way, each tree makes a categorical prediction of the predictand outcome, and probabilities are generated only in counting the proportion of trees in the forest making a particular forecast. However, this can make predictions from an individual tree very susceptible to the outcome of a particular historical case, and in some cases result in substantial overfitting. Instead, by increasing \( Z \), an RF can be allowed to make ‘impure’ leaves; at these nodes, an individual tree makes a probabilistic prediction based on the proportion of remaining training examples exhibiting each event class rather than continuing to split based on the remaining training data. Making \( S \) too large, however, can result in underfitting—lumping data as indistinguishable when there are in fact underlying discernible distinctions between remaining training examples with different labels. The last parameter, denoted \( P \), is not actually an RF algorithm parameter at all. When PCA is performed, there is always a question about the number of components to retain. Though there are some heuristics (e.g. North et al. 1982), there is no definitive method to know \textit{a priori} how many retained components \( P \) will produce the most skillful forecasts (Wilks 2011). If \( P \) is too small, valuable forecast data is discarded and predictive performance consequently suffers. However, if it is too large, the retained PCs eventually become essentially just noise,
and the RF, by fitting to these predictor values in the training data, will yield an overfit model that does not generalize to unseen data. Experiments that will not be discussed herein revealed that using information gain to determine splits and letting $B=1000$ produced skill near that of an infinitely large forest, and skill was insensitive to modifications of these settings, including modest increases in the forest size beyond this point. However, the Z-S-P parameter space are explored for the models trained and those results are presented in Appendix B.

One final consideration concerns the handling of rare event scenarios. For rare event problems, one necessarily has many more examples of the common event class in comparison to the rare class, leaving the rare class somewhat underrepresented in the learning problem, and model fitting that is done with respect to the rare class is often too dependent on a small number of examples. An approach that has been applied with some success in past studies (e.g. Ahijevych et al. 2016) is to sample training data disproportionately from the rarer classes, so that the number of training example associated with each event class are approximately equal. A comparison between this so called “balanced” sampling and unmodified “unbalanced” sampling is also made and the results presented in Appendix B.

**b. Logistic Regression**

One sensitivity experiment compares model performance as a function of the model algorithm by comparing skill of forecasts produced by RFs with those produced with logistic regression (LR). LR is in many senses a simpler model than an RF, since the structural form of the relationship between the predictors and the predictand is predefined before training. RFs, in contrast, make few assumptions about the relationships between the predictors and the predictand, allowing more diverse diagnoses of underlying relationships. However, this lack of assumptions can result in overfitting. As an application of the generalized linear model, LR assumes a linear predictors-
predictand relationship via the logit function. In LR, a single regression equation, or K equations for a multiclass problem with K categories, is computed to represent the probability of the outcome being category $k$ given the set of input predictors $x$. In particular, verifying probabilities are computed using the softmax function:

$$P(y = k | x) = \frac{e^{x^T w_k}}{\sum_{j=1}^{K} e^{x^T w_j}} \quad (A3)$$

In training a LR model, the goal is to determine the optimal weights $w_k$ associated with each predictor in order to yield the most accurate predictions for each event class. As with RF models, LR can be prone to overfitting if unconstrained. For RFs, one aforementioned approach to alleviate this problem is to increase the above-mentioned Z parameter, which stops node splitting earlier on and makes the model less tailored to the specific training data supplied to it. Complexity in LR can be thought of as being analogously represented by large weights, or regression coefficients. In order to ensure better generalizability of the trained regression equations, it is often good practice to penalize large weights through a process known as regularization. When this is done, the computation of optimal weights can be represented as a minimization problem with two terms. For 1) a matrix $Y$ with binary elements that are non-zero if and only if training example $i$ has associated verifying category $k$ and 2) a model outputting a probability matrix $P$ for each training example and category, the multinomial loss $J$ to be minimized can be computed as:

$$J(Y, P(w)) = \frac{1}{2} w^T w - \frac{1}{CN} \sum_{i=1}^{N} \sum_{k=1}^{K} Y_{i,k} \log(P_{i,k}) \quad (A4)$$

where $C$ represents the extent of regularization, with smaller values indicating that large weights are penalized more than with larger values of $C$. Alternative approaches to regularization exist (e.g. Pedregosa et al. 2011; Murphy 2012), and are explored to some degree in sensitivity experiments of Appendix B.
c. Computational Considerations

Other machine learning algorithms do not scale well to the high dimensionality of the forecast problem explored here. While time to train a model is not of primary concern for operational forecasting since it is performed only once (or periodically) offline, there are nevertheless some practical considerations; models that take months or longer to train would be unlikely to be realistic choices, for example. The “online” forecasting component, that is, the time required to take a new forecast, input it into a trained model and receive a forecast, is of operational concern, but all of the forecast techniques considered here can produce forecasts in a matter of minutes, and the small differences are not considered to be of practical concern. Using the random forest classification heuristic of considering the square root of the total number of features at each node split (Geurts et al. 2006), the computational complexity of training an RF of size $B$ from $N$ training examples with $F$ features ($N > F$) may be expressed as $O(B\sqrt{FN \log(N)})$, and may be readily parallelized across trees or within trees. Some algorithms are quadratic or even cubic (e.g. Cortes and Vapnik 1995) in the number of training examples, and do not parallelize as readily. LR is linear in the number of training examples, but requires matrix multiplication, a process that yields a computational complexity of $O(NF^2)$. PCA pre-processing, and dimensionality reduction more generally, acts both to make learning algorithms more computationally tractable and also reducing overfitting by alleviating the so-called “curse of dimensionality”.

APPENDIX B

Results: Parameter Tuning

RF model parameters were tuned for each region and lead time separately through the 4-fold cross-validation procedure employed throughout the study. Overall, the optimal parameters were
found not to vary with the two different lead times, but did vary for two of the parameters as a function of forecast region, at least to an extent; the full results appear in Table B1. For the S parameter—the number of predictors considered for each node split, the default heuristic of the square root of the total number of features was found to maximize RPSS for all regions and lead times. In all instances where both were tested, unbalanced sampling from the event classes in proportion to their true observed frequencies outperformed balanced equal sampling from each event class, in contrast to Ahijevych et al. (2016) and others; the finding appeared to be attributable to biased probabilities produced from the balanced sampling technique. For the Z parameter, the minimum number of remaining training examples in an impure parameter subspace required to perform a further node split, was generally found to be around 120. Lesser values maximized skill in the western regions, with values of 30 maximizing skill in the SW and ROCK regions, and Z=4 producing the best skill over PCST. A couple of the larger regions of the east, SE and MDWST, maximized RPSS with a value of 240, although the sensitivity between Z=120 and Z=240 was small for all regions. For P in the CTL-PCA models, skill was generally maximized with P=30, that is, retaining the 30 PCs which explain the most variance of the entire GEFS/R predictor set. For most regions, there was very limited sensitivity in the P=30–40 interval—although there was larger sensitivity outside this interval—and P=40 was found to produce slightly better skill in the NGP region. The PCST region was again the main exception, where P=60 was found to maximize cross-validation RPSS.

LR model parameters were tuned using an identical framework to ascertain the type of regularization, either based on a L1 norm which penalizes non-zero weights, or L2 norm—described in Appendix A—which penalizes large magnitude weights. L2 regularization was consistently found to produce superior results, perhaps because the number of retained PCs was already taken from the P parameter in the RF experiments, acting to nullify many potential non-zero weights of higher
numbered PCs. Unlike the RF experiments, there were occasionally some large differences in the obtained optimal regularization parameter value C between lead times within the same region. Generally, models performed better with more regularized solutions, but there were some notable exceptions, with the Day 2 NGP model and Day 3 NE model obtaining optimal C parameter values on the other end of the spectrum.

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LIST OF TABLES

Table 1. Summary of dynamical model fields examined in this study, including the abbreviated symbol to which each variable is referred throughout the paper, a description of each variable, the predictor group with which the field is associated in the manuscript text, and the highest resolution for which the field can be obtained from the GEFS/R.

Table 2. List of background predictors used in this study, and their associated symbols and descriptions.

Table 3. Summary of the models trained in this study, and the corresponding names designated to the models. ‘X’ indicates the process is performed or the information is used; a lack of one indicates the opposite. MEDIAN corresponds to the ensemble median, CTRL corresponds to the ensemble control member’s fields, and CNFDB uses the median in addition to the second-from-lowest and second-from-highest member values for each field. Horizontal radius is listed in grid boxes from forecast point; timestep denotes the number of hours between GEFS/R forecast field predictors. Slashes indicate the first number applies to the Day 2 version of the model, while the latter number applies to the Day 3 version. Letters enclosed by parentheses indicate sub-versions of models, with one parameter changed to the value adjacent to the letter. Asterisks indicate a model applies only to Day 2, and not Day 3. Otherwise, models apply to all eight forecast regions and have both Day 2 and Day 3 versions. Those models with bolded names are incorporated into the weighted blend of the final model configuration.

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Table B2. Optimal LR parameters obtained in cross-validation for the C parameter and regularization type for all lead times and regions. Evaluated for C were 0.0001, 0.0008, 0.0060, 0.0464, 0.359, 2.78, 21.54, 167.8, 1291, and 10000.
<table>
<thead>
<tr>
<th>Symbol</th>
<th>Description</th>
<th>Predictor Group</th>
<th>Grid</th>
</tr>
</thead>
<tbody>
<tr>
<td>APCP</td>
<td>Precipitation accumulation in past (3) 6 hours</td>
<td>Core</td>
<td>Native Gaussian</td>
</tr>
<tr>
<td>CAPE</td>
<td>Surface-based convective available potential energy</td>
<td>Core</td>
<td>Native Gaussian</td>
</tr>
<tr>
<td>CIN</td>
<td>Surface-based convective inhibition</td>
<td>Core</td>
<td>Native Gaussian</td>
</tr>
<tr>
<td>MSLP</td>
<td>Mean sea level pressure</td>
<td>Core</td>
<td>Native Gaussian</td>
</tr>
<tr>
<td>PWAT</td>
<td>Total precipitable water</td>
<td>Core</td>
<td>Native Gaussian</td>
</tr>
<tr>
<td>Q2M</td>
<td>Specific humidity two meters above ground</td>
<td>Core</td>
<td>Native Gaussian</td>
</tr>
<tr>
<td>T2M</td>
<td>Air temperature two meters above ground</td>
<td>Core</td>
<td>Native Gaussian</td>
</tr>
<tr>
<td>U10</td>
<td>Zonal-component of 10-meter wind</td>
<td>Core</td>
<td>Native Gaussian</td>
</tr>
<tr>
<td>V10</td>
<td>Meridional-component of 10-meter wind</td>
<td>Core</td>
<td>Native Gaussian</td>
</tr>
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<td>Q300</td>
<td>Specific humidity at 300 hPa</td>
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<td>1° × 1°</td>
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<td>Q500</td>
<td>Specific humidity at 500 hPa</td>
<td>Upper-Air Core</td>
<td>1° × 1°</td>
</tr>
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<td>Q700</td>
<td>Specific humidity at 700 hPa</td>
<td>Upper-Air Extra</td>
<td>1° × 1°</td>
</tr>
<tr>
<td>Q850</td>
<td>Specific humidity at 850 hPa</td>
<td>Upper-Air Core</td>
<td>1° × 1°</td>
</tr>
<tr>
<td>T250</td>
<td>Temperature at 250 hPa</td>
<td>Upper-Air Extra</td>
<td>1° × 1°</td>
</tr>
<tr>
<td>T500</td>
<td>Temperature at 500 hPa</td>
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<td>1° × 1°</td>
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<td>T700</td>
<td>Temperature at 700 hPa</td>
<td>Upper-Air Extra</td>
<td>1° × 1°</td>
</tr>
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<td>T850</td>
<td>Temperature at 850 hPa</td>
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<td>1° × 1°</td>
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<td>Zonal-component of 250 hPa wind</td>
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<td>Zonal-component of 500 hPa wind</td>
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<td>Zonal-component of 700 hPa wind</td>
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<td>V250</td>
<td>Meridional-component of 250 hPa wind</td>
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<td>V500</td>
<td>Meridional-component of 500 hPa wind</td>
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<td>Meridional-component of 700 hPa wind</td>
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<td>Meridional-component of 850 hPa wind</td>
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<tr>
<td>W850</td>
<td>Vertical velocity (omega) at 850 hPa</td>
<td>Upper-Air Core</td>
<td>1° × 1°</td>
</tr>
</tbody>
</table>

**Table 1.** Summary of dynamical model fields examined in this study, including the abbreviated symbol to which each variable is referred throughout the paper, a description of each variable, the predictor group with which the field is associated in the manuscript text, and the highest resolution for which the field can be obtained from the GEFS/R.
<table>
<thead>
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<th>Symbol</th>
<th>Description</th>
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<td>Median of 1-year ARIs whose closest GEFS/R grid point is the forecast point.</td>
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<tr>
<td>ARI1_LOCAL_MIN</td>
<td>Minimum of 1-year ARIs whose closest GEFS/R grid point is the forecast point.</td>
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<tr>
<td>ARI1_LOCAL_MAX</td>
<td>Maximum of 1-year ARIs whose closest GEFS/R grid point is the forecast point.</td>
</tr>
<tr>
<td>ARI10_LOCAL_MEDIAN</td>
<td>Median of 10-year ARIs whose closest GEFS/R grid point is the forecast point.</td>
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<td>ARI10_LOCAL_MIN</td>
<td>Minimum of 10-year ARIs whose closest GEFS/R grid point is the forecast point.</td>
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<tr>
<td>ARI10_LOCAL_MAX</td>
<td>Maximum of 10-year ARIs whose closest GEFS/R grid point is the forecast point.</td>
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<td>ARI1_REGIONAL_MIN</td>
<td>Minimum of 1-year ARIs that lie within the domain from which model predictors are drawn.</td>
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<td>ARI1_REGIONAL_MAX</td>
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<td>Minimum of 10-year ARIs that lie within the domain from which model predictors are drawn.</td>
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<td>ARI10_REGIONAL_MAX</td>
<td>Maximum of 10-year ARIs that lie within the domain from which model predictors are drawn.</td>
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<td>Latitude of forecast point.</td>
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<td>LON</td>
<td>Longitude of forecast point.</td>
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**Table 2.** List of background predictors used in this study, and their associated symbols and descriptions.
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<th>UAC_PCA</th>
<th>UAF_PCA</th>
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<th>CORE_CTRL</th>
<th>CORE_LSPC</th>
<th>CORE_LTME</th>
<th>CTL_LR</th>
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<td>12 (a), 6 (b*)</td>
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</table>

**TABLE 3. Summary of the models trained in this study, and the corresponding names designated to the models.**

‘X’ indicates the process is performed or the information is used; a lack of one indicates the opposite. MEDIAN corresponds to the ensemble median, CTRL corresponds to the ensemble control member’s fields, and CNFDB uses the median in addition to the second-from-lowest and second-from-highest member values for each field. Horizontal radius is listed in grid boxes from forecast point; timestep denotes the number of hours between GEFS/R forecast field predictors. Slashes indicate the first number applies to the Day 2 version of the model, while the latter number applies to the Day 3 version. Letters enclosed by parentheses indicate sub-versions of models, with one parameter changed to the value adjacent to the letter. Asterisks indicate a model applies only to Day 2, and not Day 3. Otherwise, models apply to all eight forecast regions and have both Day 2 and Day 3 versions. Those models with bolded names are incorporated into the weighted blend of the final model configuration.
<table>
<thead>
<tr>
<th>Region</th>
<th>S Parameter</th>
<th>Z Parameter</th>
<th>P Parameter</th>
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<td>SQRT</td>
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</tr>
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<td>MDWST</td>
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<td>SQRT</td>
<td>240</td>
<td>30</td>
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Table B1. Optimal RF parameters obtained in cross-validation for the Z-S-P parameter space. SQRT indicates the square root of the total number of predictors; symbols are otherwise as described in the manuscript text. Evaluated values were 1, 2, 4, 8, 16, 30, 60, 120, 240, 480 for Z, and 20, 25, 30, 40, 50, 60, 70, 80, 90, 100 for P.
### Table B2

Optimal LR parameters obtained in cross-validation for the C parameter and regularization type for all lead times and regions. Evaluated for C were 0.0001, 0.0008, 0.0060, 0.0464, 0.359, 2.78, 21.54, 167.8, 1291, and 10000.

<table>
<thead>
<tr>
<th>Region</th>
<th>Regularization</th>
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<td>MDWST</td>
<td>L2</td>
<td>0.0001</td>
<td>0.0464</td>
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<td>NE</td>
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<td>PCST</td>
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<td>0.0001</td>
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<tr>
<td>SW</td>
<td>L2</td>
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<td>L2</td>
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<td>0.0008</td>
</tr>
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