

# **Light curve completion and forecasting using fast and scalable Gaussian processes (MuyGPs)**

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## **ABSTRACT**

Temporal variations of apparent magnitude, called light curves, are observational statistics of interest captured by telescopes over long periods of time. Light curves afford the exploration of Space Domain Awareness (SDA) objectives such as object identification or pose estimation as latent variable inference problems. Ground-based observations from commercial off the shelf (COTS) cameras remain inexpensive compared to higher precision instruments, however, limited sensor availability combined with noisier observations can produce gappy time-series data that can be difficult to model. These external factors confound the automated exploitation of light curves, which makes light curve prediction and extrapolation a crucial problem for applications. Traditionally, image or time-series completion problems have been approached with diffusion-based or exemplar-based methods. More recently, Deep Neural Networks (DNNs) have become the tool of choice due to their empirical success at learning complex nonlinear embeddings. However, DNNs often require large training data that are not necessarily available when looking at unique features of a light curve of a single satellite.

In this paper, we present a novel approach to predicting missing and future data points of light curves using Gaussian Processes (GPs). GPs are non-linear probabilistic models that infer posterior distributions over functions and naturally quantify uncertainty. However, the cubic scaling of GP inference and training is a major barrier to their adoption in applications. In particular, a single light curve can feature hundreds of thousands of observations, which is well beyond the practical realization limits of a conventional GP on a single machine. Consequently, we employ MuyGPs, a scalable framework for hyperparameter estimation of GP models that uses nearest neighbors sparsification and local cross-validation. MuyGPs allows us to quickly train models on light curve data in seconds on a single workstation. In this manuscript we explore light curve completion and forecasting, and compare embeddings of the light curve data into multi-dimensional spaces to take advantage of daily and yearly periodicity. We show that our method outperforms feedforward DNNs both in terms of accuracy and quantity of training data required.

## 1. INTRODUCTION

Photometric light curves are a series of observations that track the brightness of an object over a period of time. They can be used to characterize the dynamic properties of a system. Astronomers frequently analyze light curves to understand a whole range of astrophysical topics including: detailed physics within a star [6], the discovery of exoplanets [2], classifying distant supernovae [23], and characterizing the population of near earth asteroids [33].

Space Domain Awareness (SDA) involves monitoring, detecting and understanding the population of earth-orbiting bodies or resident space objects (RSOs). SDA is of increasing importance due to the incipient growth in the number of space objects and debris orbiting the earth, due in large part to the recent commercial development of satellites. The number of such RSOs is expected to grow by orders of magnitude in the next decade. While SDA is concerned with maintaining custody of orbiting objects, it also prioritizes the rapid identification of changes to and anomalous behavior of the many varying orbital systems.

Fortunately, light curves are valuable for inspecting RSOs in much the same way as astrophysical phenomena, and can enable critical information for SDA. The brightness of an RSO depends on structural features like size, shape, and material composition as well as the geometry between the object, the sun and observer. The proliferation of low cost commercial-off-the-shelf (COTS) ground-based telescopes has made the generation of light curves from RSOs easier to produce. Furthermore, many constellations of ground-based telescopes have been tasked with tracking RSOs for SDA. These factors have enabled the relatively cheap production of large volumes of RSO light curves, which has prompted practitioners to apply automation to analyze them at scale.

Practitioners have recently applied machine learning to light curves of RSOs to solve various SDA tasks. For example, comparing the light curve of an unknown RSO to a catalog of known RSOs is useful for predicting features like shape [21, 13], material composition [9], and general categories or classes [20, 17, 12]. Furthermore, forecasting RSO light curves into the future is useful for detecting deviations from the expected patterns-of-life in near-real-time, affording the detection of anomalous events such as maneuvers [30, 11] or configuration changes.

The goal of machine learning in this context is to learn a function mapping an input space (typically the time domain) to an observation space, e.g. magnitude, based upon independent and identically distributed (i.i.d.) samples from a distribution of input-observation pairs. We will refer to this data distribution as the “target distribution”. A machine learning model is successful when it is able to accurately predict the observation of an unseen input drawn from the target distribution.

Both diffusion-based or exemplar-based methods have been traditionally used in image or time-series completion problems. Diffusion-based methods [31] use thermal diffusion equations to propagate information from surrounding regions into missing regions, they are mostly effective when the gaps are small and tend to smooth details out for larger problems. Exemplar-based methods [7] use greedy algorithms to apply patches of training data to missing regions which can produce implausible patterns.

Deep Neural Networks (DNNs) are an especially popular tool to model light curves in the literature due to their expressiveness and generalization capabilities. A DNN is a type of representation learning model — a machine learning model that learns an appropriate feature representation of the data in addition to producing predictions. DNNs iteratively transform the input space into latent feature spaces using linear functions that are “activated” by element-wise nonlinear functions. DNNs also happen to be universal function approximators — it is possible to approximate any continuous function to an arbitrary level of precision using a sufficiently large DNN. The persistent popularity of DNNs derives from several sources, such as the widespread availability of hardware accelerators like graphical processing units (GPUs), advanced stochastic optimization techniques that aid in their training, and the development of user-friendly software libraries such as Tensorflow [1] and PyTorch [27].

Although DNNs have many positive features, they also have some drawbacks that are particularly notable in the light curve problem. First, modern DNN architectures typically consist of a very large number of parameters that must be trained. Training such models generally consumes a large amount of computing resources, as the model and its gradient are evaluated over many iterations in order to refine parameter values. In addition to computational expense, training a large model typically requires a large amount of data. The literature has observed a roughly linear relationship between model size and the amount of labelled data required to train it [32]. This means that learning an accurate light curve model can require a large number of observations in general.

Second, while DNN models lend themselves to high dimensional feature spaces they tend to struggle with very small numbers of dimensions. Feature engineering can often solve this problem, especially in time series scenarios where some notion of a moving window usually serves as a feature space. However, this strategy is most successful when the observation cadence of the light curve is high. Upcoming ground-based surveys of deep space, such as the Legacy Survey of Space and Time (LSST) are expected to incidentally capture many RSO images from which light curves can be extracted. However, these light curves will be irregular, sparse, and have low-dimensional feature spaces, increasing the challenge of applying existing techniques.

Third, effective training typically requires a large volume of training data that is complete and representative of the target distribution i.e. a large model requires a large amount of training data that is i.i.d. according to the target distribution. In addition to informing prediction accuracy, data independence and size helps complex DNN models avoid overfitting and simply memorizing the training data. However, collections of physical measurements are often limited by the realities of sensor availability and physical obstruction. For example, weather affects optical astronomical measurements by introducing correlated noise or blocking the desired object from view entirely. Furthermore, it is in general desirable to design a model that can be alternatively applied to several different RSO inference problems, i.e. several distinct target distributions. A transfer learning approach—one where the model is at least partially trained on data from a possibly different distribution than the target distribution [32]—could address this need for a large volume of training data. Indeed, transfer learning has been applied to satellite aperture radar images [29], radiofrequency interference characterization [19], and classifying RSOs using light curves [12]. However, transfer learning ultimately relies on features learned from a potentially unrelated dataset, which may lead to inefficient or inaccurate conclusions.

Generative models such as Variational Autoencoders (VAEs) and Generative Adversarial Networks (GANs) are additional alternatives from the machine learning literature that address data efficiency. VAEs are autoencoders that impose structure upon their learned encoder and decoder functions to ensure that the latent representation of the data encodes the target distribution. Researchers have successfully applied VAEs to learn shapes from the light curves of RSOs [13]. GANs conversely simultaneously train models that generate samples meant to mimic the target distribution while distinguishing between real and synthetic samples. Practitioners have recently used GANs to classify stars based upon their light curves [14]. However, GANs can be challenging and expensive to train. GANs infer a distribution from a very small training dataset and can suffer from mode collapse, non-convergence or general instability.

This brings us to the fourth point — DNNs have difficulty quantifying the uncertainty in their predictions. It is generally difficult to determine when a DNN is extrapolating — predicting the response on data from a different distribution or in a different region of the training data. This is problematic in scientific applications because overconfident predictions can lead to incorrect inference and unnatural results that are difficult to interpret. Furthermore, it is important for decision makers to distinguish between low and high-confidence predictions when drawing conclusions for SDA.

Although uncertainty quantification methods are in research, most practical models employed by practitioners provide only point predictions with no obvious mechanism to measure model confidence. Recent attempts to provide uncertainty quantification to DNNs attempt to learn prediction intervals (i.e. error bars) in addition to point predictions, but this literature is still developing and there is as yet no consensus solution [18]. Others attempt a hybrid approach where a Gaussian process is appended to the last layer of a DNN [4].

In this manuscript we propose Gaussian Process (GP) models as alternatives for the light curve modeling problem. Like DNNs, GPs are representation learning methods that are data-driven universal function approximators. A GP is a type of kernel method that implicitly and non-linearly maps input features into a possibly infinitely dimensional inner product space. Kernel methods use a parameterized kernel function to cheaply compute inner products in this implicit space to predict unknown responses. GPs can also be thought of as the generalization of a multivariate normal distribution, where all data within a defined domain is assumed to be jointly normally distributed with the covariance defined by the kernel function. Therefore, predictions from GP models are probabilistically defined through conditional probability. GPs are attractive for scientific applications due to this inherently Bayesian inference model, which produces explicit uncertainty quantification (Gaussian distributions) of its predictions. Furthermore, GPs have been shown to outperform DNNs in data-starved regimes [26] and are ideal models for low-dimensional feature spaces [25].

However, conventional GPs have very poor scaling in the number of training observations, which has previously limited their application to the light curve modeling problem. Both realizing the predictions from the GP model and evaluating the likelihood function in training require cubic computation in the number of data points, and require quadratic memory to store the kernel matrix. While this practical drawback has tended to limit the application of GPs

to small data problems, scalable approximate GP methods have proliferated in the literature [16, 22]. These methods typical tradeoff accuracy for speed. However, the approximate GP method MuiGPs offers a selection of settings that demonstrates superior accuracy or computational scaling in several datasets [25, 26, 5].

Therefore, in this paper we use the MuiGPs approximate GP estimation method [25]. MuiGPs uses nearest neighbors sparsification and a batched leave-one-out cross-validation objective function to train a GP-like model in nearly linear time in the number of data observations. Investigators have successfully applied MuiGPs to cosmology image processing problems [15, 26, 5].

In this paper, we propose a method of light curve interpolation and prediction using MuiGPs that gives both predictions and uncertainty quantification of those predictions. We use this method in two modes:

1. Interpolation, the prediction of unobserved magnitudes in the past, and
2. Forecasting, the prediction of unobserved future magnitudes.

Interpolation is useful as a data preprocessing utility for catalogs. Interpolation allows us to fill in magnitude observations for sparse, irregular light curves to make them suitable for downstream comparison tasks such as shape, pose, size, type, or material estimation. Forecasting is useful to detect deviations from expected future patterns-of-life that correspond to anomalies. The posterior variance provided by GPs is especially useful to determine what constitutes a significant deviation from expected behavior.

In section 2, we provide background on the light curve data itself, data processing, as well as the machine learning methods we utilize in our comparative study. Then in section 3 we describe several numerical studies that demonstrate the comparative performance of several choices of our GP method to that of a DNN and how the uncertainty quantification provided by our method can be used to detect anomalies. Finally, in section 4, we discuss our conclusions, the importance of our findings, limitations, and future work associated with our methodology.

## 2. METHODOLOGY

### 2.1 Light curves definition

Light curves are time series of the brightness of resident space objects over long periods. They are obtained by observing the same object, night after night, for several days or even years. One way to visualize them is in two dimensions with one axis representing the time of day (or solar phase angle) and the other axis representing the days while the pseudo-color indicates the brightness, as in Fig. 1.

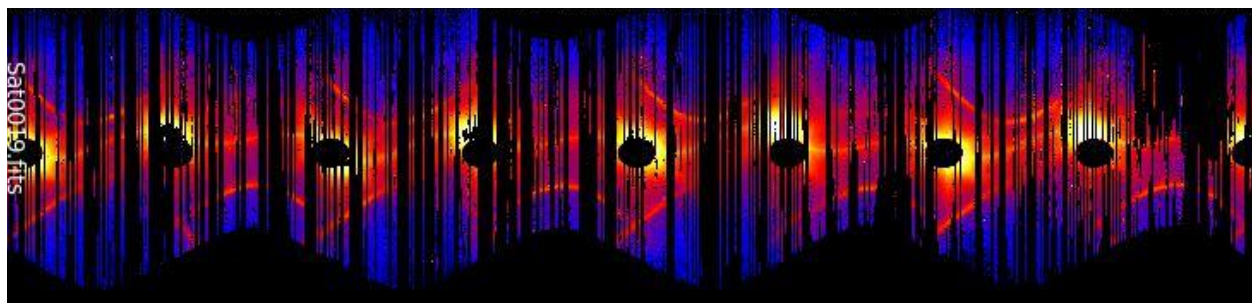


Fig. 1: The light curve of a single object over 4 years. Each column represents a single day. The color scale represents the brightness. The black stripes and *holes* highlight the sparsity of observations; the former are due to weather, the latter to eclipse. The seasonality and yearly periodicity induces the outer shape. Courtesy of [24].

Light curves of RSOs contain potentially detailed information. For example, one could detect dust accumulating on highly reflective solar panels and deduce the age and life expectancy of a satellite. Higher intensity reflections also have the potential to inform about the shape of the observed object since different facets of a multi-faceted object would reflect differently. In addition, satellites can maneuver or rotate, in which case, the reflected light will deviate from the expected pattern. Those deviations can be detected and increase SDA knowledge.

Observations can be limited by the time of day, weather, sensor saturation, and eclipses (see Fig. 2). Data is only available at night, and the measurements are less accurate near dusk and dawn. The weather can preclude observations for hours and sometimes days at a time. When the reflected light from the sun is particularly bright, sensors can saturate leading to missing data, though this particular case is usually fairly obvious when looking at the brightness of surrounding available data. Lastly, periodically, the light from a satellite will be eclipsed by the earth.

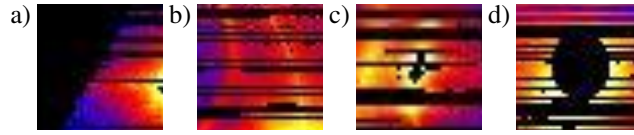


Fig. 2: Missing data due to a) time of day, b) weather, c) sensor saturation, and d) eclipse.

To test our codes we use a light curve dataset provided by Dave Monet [24] of 43 satellites. All satellites in the dataset are from the public catalog ([www.space-track.org](http://www.space-track.org)). We selected 13 distinct RSOs, those manually flagged by Dave Monet as *nominal*, for our analysis. There are approximately 500,000 data points per object. All observations were taken from a single camera in Flagstaff AZ between September 2014 and September 2018. The dataset contains the brightness (magnitude; band is approximately Johnson V) as well as the measurement error in brightness (uncertainty).

## 2.2 Processing

To test our prediction capabilities, we select a portion of each light curve as test data. We use time periods ranging from a few hours (3 hours) to multiple weeks (2 weeks). And we either select the test data at a random time within the time series (interpolation) or at the end (extrapolation). To guarantee that the selected period does not fall at a time lacking observations (e.g. during the day or during a cloudy night) we reject those that have fewer than 90% of the data we'd expect to find in that time interval if the data was uniformly distributed. In addition, we apply that same condition to the preceding interval to ensure that we're not accidentally interpolating for longer that we would expect.

Since the light curves have some daily and yearly periodicity, we compare several embeddings of the light curve data into multi-dimensional spaces. The reference is the 1D model which is just the original time series. The 2D model has one dimension representing times of day as real numbers in the  $[0, 1]$  interval, and another dimension representing days as integers, starting at zero on the first day of observation. The 3D model is like the 2D model but with an additional dimension for years, as integers between 0 and 4, and with a modified days dimension modulo 365. Note that all the input dimensions are eventually normalized to the interval  $[0, 1]$  as is customary.

An alternate 2D model with year and time of year as opposed to day and time of day was considered but dismissed for simplicity and for not being as intuitive and interesting. We observed that day-to-day correlation is stronger than year-to-year in our dataset.

Note that it is more traditional to model this type of data periodicity using a periodic kernel, but the MuiGPs estimation framework depends on a kernel sparsity that is ultimately not validated with that kernel type. Therefore, these embeddings are a novel way to replace such a periodic kernel while maintaining the computationally efficient framework.

## 2.3 Gaussian processes

We will consider throughout a light curve to be a univariate response  $Y : \mathcal{X} \rightarrow \mathbb{R}$ , where  $\mathcal{X} \subseteq \mathbb{R}^d$  is the observation space along  $d$  time dimensions. In preprocessing we de-trend the data, and so  $Y$  therefore has zero mean. In modeling  $Y$  with a GP, we assume that it is drawn from a Gaussian process distribution. This means that  $Y$  follows a multivariate Gaussian distribution at any finite set of  $n$  points  $X = (\mathbf{x}_1, \dots, \mathbf{x}_n) \in \mathcal{X}^n$ . However, in reality measurement noise perturbs our observed values of  $Y$  at locations  $X$ . We assume that each measurement is perturbed by homoscedastic noise that are i.i.d.  $\mathcal{N}(0, \epsilon)$ . That is,

$$Y(X) = (Y(\mathbf{x}_1), \dots, Y(\mathbf{x}_n))^T \sim \mathcal{N}(\tilde{0}, Q_\theta(X, X)), \quad (1)$$

$$Q_\theta(X, X) = \sigma^2 (K_\theta(X, X) + \epsilon I_n), \quad (2)$$

where  $\mathcal{N}$  is the multivariate Gaussian distribution,  $\tilde{0}$  is the  $n$ -dimensional zero vector,  $\sigma^2$  is a variance scaling term,  $I_n$  is the  $n \times n$  identity matrix, and  $K_\theta(X, X)$  is an  $n \times n$  positive definite, symmetric covariance matrix between the

elements of  $X$  that is controlled non-linearly through kernel function  $K_\theta(\cdot, \cdot)$  with hyperparameters  $\theta$ .  $Q_\theta(X, X)$  is  $K_\theta(X, X)$  that is perturbed by the measurement noise prior  $\varepsilon$  and scaling parameter  $\sigma^2$ .

Similarly, any additional (possibly unobserved) datum  $\mathbf{x}^* \in \mathcal{X}$  is also jointly normal with observed data  $X$  by the GP assumption. Thus, the conditional distribution for the response of  $Y$  at  $\mathbf{x}^*$  given responses observed at  $X$  and  $\theta$  is also multivariate normal with mean and variance

$$Y_\theta(\mathbf{x}^* | X) = K_\theta(\mathbf{x}^*, X) Q_\theta(X, X)^{-1} Y(X), \text{ and} \quad (3)$$

$$\text{Var}(Y_\theta(\mathbf{x}^* | X)) = K_\theta(\mathbf{x}^*, \mathbf{x}^*) - K_\theta(\mathbf{x}^*, X) Q_\theta(X, X)^{-1} K_\theta(X, \mathbf{x}^*), \quad (4)$$

where  $K_\theta(\mathbf{x}^*, X) = K_\theta(X, \mathbf{x}^*)^\top$  is the cross-covariance matrix between  $\mathbf{x}^*$  and the elements of  $X$ .

GPs are typically trained by maximizing the log-likelihood of the observations  $Y(X)$  with respect to  $\theta$ . This log-likelihood function possesses the following form:

$$\log(L(\theta, Y(X))) = -\frac{p}{2} \log(2\pi) - \frac{1}{2} \log(|Q_\theta(X, X)|) - \frac{1}{2} Y(X)^T Q_\theta(X, X)^{-1} Y(X). \quad (5)$$

However, evaluating Equation 5 is  $O(n^3)$  computation and  $O(n^2)$  in memory, which is intractable for all but relatively small data. A MuiGPs model rewrites equations 3 and 4 as

$$\hat{Y}_\theta(\mathbf{x}^* | X_{N^*}) = K_\theta(\mathbf{x}^*, X_{N^*}) Q_\theta(X_{N^*}, X_{N^*})^{-1} Y(X_{N^*}), \text{ and} \quad (6)$$

$$\text{Var}(\hat{Y}_\theta(\mathbf{x}^* | X_{N^*})) = K_\theta(\mathbf{x}^*, \mathbf{x}^*) - K_\theta(\mathbf{x}^*, X_{N^*}) Q_\theta(X_{N^*}, X_{N^*})^{-1} K_\theta(X_{N^*}, \mathbf{x}^*), \quad (7)$$

where  $X_{N^*}$  are the nearest neighbors of  $\mathbf{x}^*$  in  $X \setminus \{\mathbf{x}^*\}$ . MuiGPs trains  $\theta$  in terms of some loss function  $\ell(\cdot, \cdot)$  over a sampled batch  $B = (\mathbf{x}_1^*, \dots, \mathbf{x}_b^*) \subseteq X$  of size  $b$  by minimizing an objective function  $Q(\theta)$ . Minimizing  $Q(\theta)$  allows us to train  $\theta$  without evaluating the expensive determinant in the GP likelihood. When we set  $Q(\theta)$  to leave-one-out cross-validation and  $\ell_\theta$  to mean squared error, training  $\theta$  reduces to minimizing the function

$$Q(\theta) = \frac{1}{b} \sum_{i \in B} \left( Y(\mathbf{x}_i^*) - \hat{Y}_\theta(\mathbf{x}_i^* | X_{N_i^*}) \right)^2. \quad (8)$$

Note other loss functions can be employed in this framework, but this mean squared error function has been demonstrated as performant [25]. We use Bayesian optimization to optimize Equation 8 to train  $\theta$  in our experiments.

## 2.4 A standard neural network model for benchmarking

Neural networks have achieved state-of-the-art results on common benchmark problems and are now ubiquitous in scientific applications. Although the focal point of this work is the application of Gaussian processes to light curve modeling, we benchmark the predictive performance of MuiGPs against a standard deep neural network model to ensure its predictive accuracy. We select a fully-connected architecture, as the structure of the feature space is not suited to convolutional models. Moreover, recurrent neural networks, including LSTMs, are another popular class of models that are useful in forecasting settings, but not interpolation problems.

Fully-connected neural networks can be represented as the composition of a series of affine transformations and nonlinear activation functions. The composition of an individual affine transformation and nonlinear activation function constitutes a layer of the fully-connected network. Each affine transformation in the network is parameterized by weights  $\mathbf{W}_i \in \mathbb{R}^{n_i \times n_{i+1}}$ , where  $n_i$  is the dimension of the input to the  $i^{\text{th}}$  layer and  $n_{i+1}$  is the output dimension, and a bias vector  $\mathbf{b}_i \in \mathbb{R}^{n_{i+1}}$ . We label the activation function in the  $i^{\text{th}}$  layer  $\sigma_i$ . Let  $\mathcal{F}_{NN}(\mathbf{x}) : \mathbb{R}^{n_{in}} \rightarrow \mathbb{R}^{n_{out}}$  be a neural network. Then,

$$\mathcal{F}_{NN}(\mathbf{x}) = \sigma_n(\mathbf{W}_n \sigma_{n-1}(\mathbf{W}_{n-1} \sigma_{n-2}(\mathbf{W}_{n-2}(\dots) + \mathbf{b}_{n-2}) + \mathbf{b}_{n-1}) + \mathbf{b}_n). \quad (9)$$

We train a fully-connected neural network using the day of the year and time of day as the independent variables (features), with the target set to the magnitude (normalized to take on values between 0 and 1 by dividing by the largest magnitude observed). The model architecture features ReLU activation functions and 5 layers of sizes  $2 \times 200$ ,  $200 \times 200$ ,  $200 \times 100$ ,  $100 \times 20$ , and  $20 \times 1$ . There are 62,420 total parameters in the model, roughly  $1/8 - 1/5$  the number of training samples depending on the test case. We train the neural network in batches of size 128 using the Adam optimizer for 100 epochs. We use the mean-squared error loss function with 20 percent of the training data

withheld for validation, the training rate set to  $10^{-3}$ , and a learning rate exponential decay factor of 0.5 applied every 30 training epochs. We set the ratio of the number of parameters in the model to the number of data points on the order of  $1/10$ , a commonly accepted ratio in the deep learning community. We selected these training hyperparameters based on common values used in deep learning to achieve the best combination of training loss decay and validation error minimization.

### 3. RESULTS

First we compare the predicting power of Gaussian processes for the different embeddings of the input data described in subsection 2.2. For each of the 13 light curves in our dataset, we randomly select a total of 20 time intervals, 5 for each time duration in  $\{3 \text{ hours}, 1 \text{ day}, 1 \text{ week}, 2 \text{ weeks}\}$ , and use those intervals as test data while using the rest as training data. The accuracy of the prediction is measured as the Root Mean Square Error (RMSE) divided by the extent (max - min) of the magnitude in the test data.

We use the common radial basis function (or RBF) to define our kernel in all of our experiments. This means that we use the kernel function

$$K_\ell(x, x') = \exp\left(-\frac{\|x - x'\|_2^2}{2\ell^2}\right). \quad (10)$$

The RBF kernel has length scale parameter  $\ell$ , and our model additionally has the measurement noise variance prior  $\varepsilon$  and variance scaling parameter  $\sigma^2$ . We fix  $\varepsilon = 10^{-5}$  in our experiments. We train  $\ell$  by optimizing equation Equation 8 via Bayesian optimization. However, no prediction-based objective function is sensitive to  $\sigma^2$ , and so we optimize it according to the analytic procedure described in Section 2 of [25].

Fig. 3 shows the distribution of the resulting comparison of the predictions to the truth for each embedding. With a median of 0.066, the 2D embedding yields more than 3 times better predictions than the original 1D embedding (median: 0.24). The better prediction is possible because the distance between similar data points across days is reduced, thanks to the extra dimension, allowing the GP to more readily use them during interpolation. However we see that adding a third dimension encoding the yearly periodicity does not yield significant improvements over the 2D embedding since the results are quasi-identical (median: 0.065). Although the distributions of 2D and 3D embeddings are similar, from this on, we focus on the 2D embedding and use it exclusively.

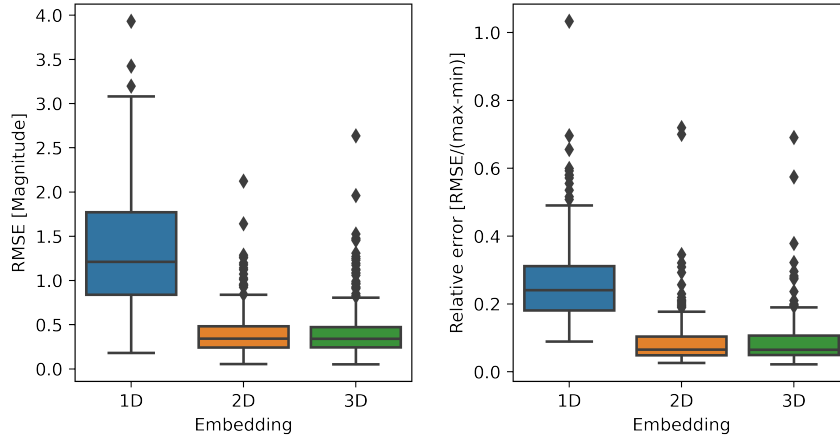


Fig. 3: Accuracy of predictions for different data embeddings as a standard box plot representing the distribution of results of completing 5 random gaps of 4 distinct durations ranging from 3 hours to 2 weeks for each of the 13 light curves in the dataset. 1D is the raw time series of magnitudes, 2D takes into account daily periodicity, and 3D takes into account both daily and yearly periodicity. While GPs benefit greatly from the daily periodicity information, the effect of adding information about yearly periodicity is comparatively negligible.

Part of the motivation for being able to predict data in light curves is to complete gaps and missing data points. A first



benchmark is to be able to complete randomly selected data points throughout a light curve. We randomly select either 5%, 10%, or 20% of the total number of data point available in each of the 13 light curves in the dataset, 5 times for each proportion. Then, for more realistic interpolation, we select consecutive data points representing gaps of either 3 hours, 1 day, 1 week, or 2 weeks, with 5 different random starting instants for each 13 light curve. Lastly we repeat the same procedure but by selecting gaps at the end of each light curve to evaluate how GPs perform for extrapolation.

Fig. 4 shows the performance of GP predictions for all of these interpolation and extrapolation tasks. For random interpolation, for all percentages (5%, 10%, or 20%), the performance is similar and very good with a median relative RMSE of about 0.03. For gaps, either interpolation or extrapolation, there is more variation, both for a given gap duration, and across durations.

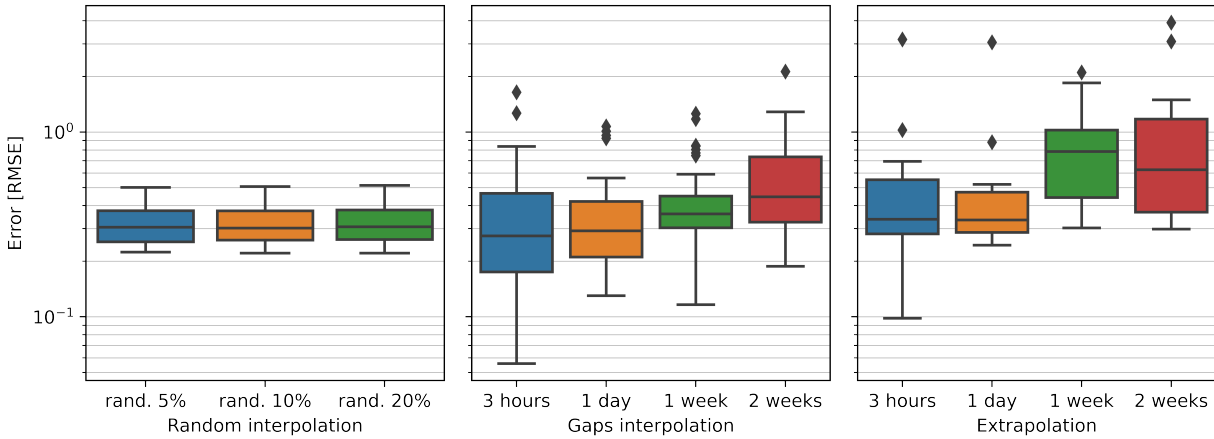


Fig. 4: Comparison of the predicting power of GPs (with 2D embedding) for various interpolation and extrapolation tasks. The left pane is for interpolating randomly selected datapoint, representing from 5% to 20% of the entire light curves. The middle pane is for interpolating gaps of durations ranging from 3 hours to 2 weeks. The right pane is for extrapolating gaps of similar durations at the end of each light curves. Random data points are interpolated with great accuracy, even when they represent a significant portion of the data available. Interpolating gaps over short periods of time is more challenging.

To compare the performance of GPs compared to DNNs, we ran the same task of predicting one day of missing data in a single 4 year-long light curve with training data exclusively from that single light curve on the same machine. As shown in Fig. 5, GPs achieve better accuracy in only a fraction of the time.

Fig. 6 illustrates a single example of the quality of the GP prediction compared to raw measurements for a day of missing data using the 3D embedding. The mean prediction follows the trend of magnitudes well as expected. As is typical in GP models, our kernel model is assumed to be stationary and homoscedastic, meaning that we learn a single set of hyperparameters that best models all the training data (across all magnitudes). In this interval pictured, we see that the data at high magnitudes seem to have more variance than the data collected at lower magnitudes. Our model is agnostic to any change in variance in the data itself, but will on average provide desired uncertainty quantification. Note this one timegap is a single interval that is in itself time-correlated in magnitude and therefore variance regime. Therefore, the coverage of the 95% interval from this one sample can be different than the desired overall level, but when we consider many independent time intervals of this type, the uncertainties will average to the desired 95% confidence level. In future work more flexible and complex GP kernel functions could be designed with inhomogeneous or non-stationary kernels to correct for this potential pattern in magnitude to improve uncertainties in each individual sample region. If the magnitude distribution in our prediction interval differs from that of our training data then there is a model discrepancy in our approach that could explain our observed variances.

But the true power of GPs comes from their ability to predict a full posterior distribution and not just a mean. The predicted covariance can be used to detect areas where the prediction differs from the measurement by a significant amount. One application of particular interest for SDA is the detection of anomalies — potential maneuvers or state changes — in light curve data. Fig. 7 shows two example of such anomalies. In both cases, two factors make the



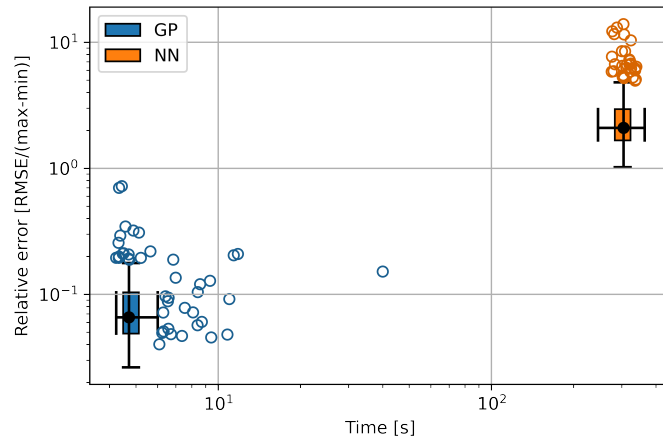


Fig. 5: Comparison of the performance of our Gaussian process method with an equivalent neural network. The accuracy is measured by the RMSE scaled by the extent of the magnitude in the test dataset. The computing time is the total of the training and prediction time as those cannot be separated for the GP, and the prediction time is negligible compared to the training time for the NN. On the same data, the Gaussian process achieves a better accuracy in a fraction of the time.

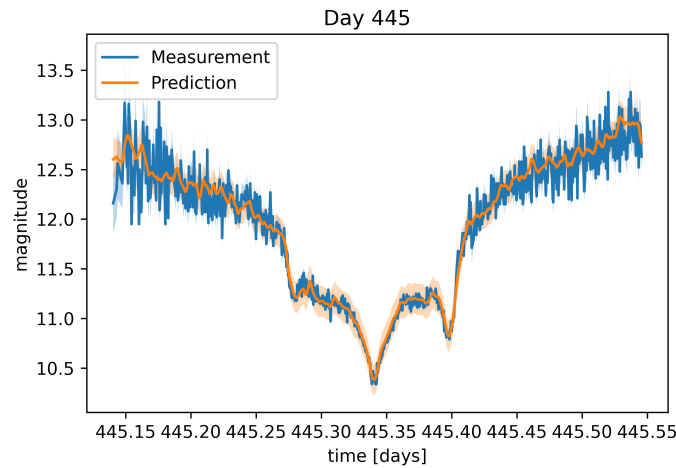


Fig. 6: GP prediction (orange) compared to measurement (blue) for 1 day of missing data. The orange shaded region corresponds to the 95% confidence interval of the prediction while the blue shaded region corresponds to the measurement standard error. This prediction was obtained using a RBF kernel with three dimensional training data. The GP does a very good job a predicting details while effectively filtering noise.

anomalies noticeable. Firstly, the prediction differs from the measurement for a small but significant period of time, and the difference exceeds two times the predicted standard deviation. And secondly, the measurements from the surrounding days look similar to those of the focal day — and to the prediction — except for that particular period. The combination of those two factors indicate that it is indeed the measurements that are anomalous and not the predictions.

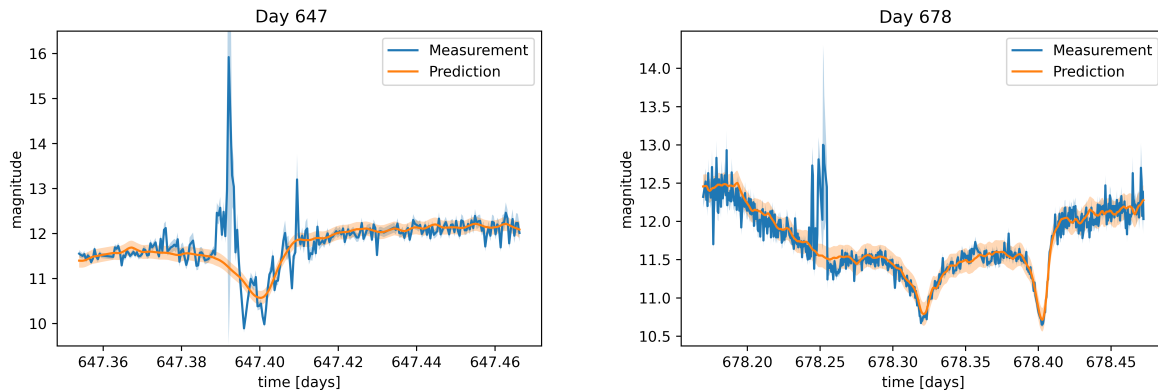


Fig. 7: Example of detected anomalies. The prediction is smooth and matches the measurement well except for a short time interval (near 647.39 and 678.25 days respectively) where the measurement suddenly exceeds its expected value based on the prediction and confirmed by measurements from the surrounding days.

#### 4. CONCLUSION

We have shown Gaussian processes (GPs) using MuiGP to be a capable and pertinent tool for analyzing light curve data. Unlike other machine learning methods that can have millions of parameters, Gaussian processes are simple in that they do not have a large architecture search and in that they only have a few parameters to be estimated. Unlike traditionally GP estimation methods, the implementation of the MuiGPs methodology allows astronomers to apply GP methods to very large data with very little computational or memory burden. In addition to their improved accessibility, we have shown that in our designed experiments on light curves, Gaussian processes are both more accurate and faster than an example neural network. Further, the uncertainty quantification we get from Gaussian processes allows us to identify statistically distinct deviations from the pattern-of-life.

As seen above, while Gaussian processes are able to ingest raw uni-dimensional time series, encoding the periodicity of the light curve data through additional dimensions yields much better predictions when we predict large gaps of observations. This is likely due to the distinct periodic trends in the light curves. However, we observed that encoding the daily periodicity was sufficient and that encoding the yearly periodicity did not further improve prediction. A first explanation is that surrounding days are typically more similar than surrounding years, so the majority of “nearest neighbors” used in the GP model would naturally be selected from surrounding days rather than from surrounding years.

We have also shown how Gaussian processes compare advantageously to neural networks both in terms of computing time during training and prediction accuracy. Note that our metric for computing time combines both training and prediction time from the two models. This metric masks the tradeoff that a trained neural network can make additional predictions efficiently whereas training a Gaussian process using MuiGPs is very fast, but the prediction time given a trained model is much slower comparatively. [10] demonstrate a way to improve prediction time of a similar kernel interpolation with an additional approximation. However, in the online scenario of the light curve problem where one may want to perform this analysis in near realtime, one would both retrain and predict simultaneously so the GP method is significantly preferred.

Both methods can be parallelized to take advantage of High Performance Computing (HPC). For DNNs, parallelizations techniques are readily available, for instance in PyTorch, with the advent of GPUs and TPUs. For MuiGPs, parallelizations efforts are in progress and will be released in a different publication. We ran all of the experiments in

this manuscript on a single core of a commodity laptop using the software library MuyGPyS [28].

Future applications will use MPI [8] and JAX [3] to scale model training and evaluation to multiple compute nodes using many CPUs and GPUs in parallel. These scalability features will enable applications that scale to the observation sizes of LSST, which are anticipated to involve hundreds of millions of space objects.

Possible future directions of research involve comparing GP training from one set of light curves and using these trained parameters to make predictions for a different set of unseen light curves. This will determine whether it is necessary to fit parameters to each light curve. Finally, since our method gives fully interpolated light curves, future work could use these de-noised full time series for further SDA studies.

In summary, we believe our GP method is a valuable method for light curve interpolation and future prediction for three scientific tasks of particular interest. First, by de-noising and interpolating missing data, GPs can be part of a pre-processing pipeline before feeding the data to further algorithms or software that cannot work with raw data. As shown in figures 1 and 2, raw light curves can have a lot of gaps frequently spanning multiple days so being able to interpolate those gaps is crucial. Second, by being able to extrapolate data over several hours, days, or even weeks, GPs can be used to predict the expected behavior of RSOs, which can serve in forecasting and collision avoidance in SDA. Last and not least, by quantifying the uncertainty of predictions, GPs allow for automatic change detection. Indeed, being able to predict a full posterior distribution enables discerning anomalies in measurements from anomalies in prediction, since uncertain predictions can easily lead to false positives. In future research, it should be possible to use tracking data and/or documented known maneuvers to fine-tune and benchmark our maneuver detection capabilities.

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