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Overview of GeoLifeCLEF 2021: Predicting species distribution from 2 million remote sensing images

Titouan Lorieul, Elijah Cole, Benjamin Deneu, Maximilien Servajean, Pierre Bonnet and Alexis Joly

Abstract
Understanding the geographic distribution of species is a key concern in conservation. By pairing species occurrences with environmental features, researchers can model the relationship between an environment and the species which may be found there. To advance research in this area, a large-scale machine learning competition called GeoLifeCLEF 2021 was organized. It relied on a dataset of 1.9 million observations from 31K species mainly of animals and plants. These observations were paired with high-resolution remote sensing imagery, land cover data, and altitude, in addition to traditional low-resolution climate and soil variables. The main goal of the challenge was to better understand how to leverage remote sensing data to predict the presence of species at a given location. This paper presents an overview of the competition, synthesizes the approaches used by the participating groups, and analyzes the main results. In particular, we highlight the ability of remote sensing imagery and convolutional neural networks to improve predictive performance, complementary to traditional approaches.

Keywords
LifeCLEF, evaluation, benchmark, biodiversity, presence-only data, environmental data, remote sensing imagery, species distribution, species distribution models

1. Introduction
In order to make informed conservation decisions, it is essential to understand where different species live. Citizen science projects now generate millions of geo-located species observations every year, covering tens of thousands of species. But how can these point observations be used to predict what species might be found at a new location?

A common approach is to build a species distribution model (SDM) [1], which uses a location’s environmental covariates (e.g. temperature, elevation, land cover) to predict whether a species is likely to be found there. Once trained, the model can be used to make predictions for any location where those covariates are available.
Developing an SDM requires a dataset where each species observation is paired with a collection of environmental covariates. However, many existing SDM datasets are both highly specialized and not readily accessible, having been assembled by scientists studying particular species or regions. In addition, the provided environmental covariates are typically coarse, with resolutions ranging from hundreds of meters to kilometers per pixel.

In this work, we present the results of the GeoLifeCLEF 2021 competition which is part of the LifeCLEF evaluation campaign [2] and co-hosted in Eighth Workshop on Fine-Grained Visual Categorization (FGVC8)\(^1\) at CVPR 2021. This competition is the fourth GeoLifeCLEF challenge. In the first two editions, GeoLifeCLEF 2018 [3] and GeoLifeCLEF 2019 [4], each observation was associated only with environmental features given as vectors or patches extracted around the observation. Like last year’s competition, GeoLifeCLEF 2020 [5], GeoLifeCLEF 2021 is aimed at bridging the previously mentioned gaps by (i) sharing a large-scale dataset of observations paired with high-resolution covariates and (ii) defining a common evaluation methodology to measure the predictive performance of models trained on this dataset. The dataset is based on over 1.9 million observations of plant and animal species. Each observation is paired with

\(^1\)https://sites.google.com/view/fgvc8
high-resolution remote sensing imagery – see Figure 1 – as well as traditional environmental covariates (e.g. climate, altitude and soil variables). To the best of our knowledge, this is the first publicly available dataset to pair remote sensing imagery with species observations. Our hope is that this analysis-ready dataset and associated evaluation methodology will (i) make SDM and related problems more accessible to machine learning researchers and (ii) facilitate novel research in large-scale, high-resolution, and remote-sensing-based species distribution modeling.

2. Dataset and evaluation protocol presentation

Data collection. The data for this year’s challenge is the same as last year reorganized in a more easy-to-use and compact format. A detailed description of the GeoLifeCLEF 2020 dataset is provided in [6]. In a nutshell, it consists of 1,921,123 observations covering 31,435 species (mainly plants and animals) distributed across US (1,097,640) and France (823,483), as shown in Figure 2. Each species observation is paired with high-resolution covariates (RGB-IR imagery, land cover and altitude) as illustrated in Figure 1. These high-resolution covariates are resampled to a spatial resolution of 1 meter per pixel and provided as $256 \times 256$ images covering a $256m \times 256m$ square centered on each observation. RGB-IR imagery come from the 2009-2011 cycle of the National Agriculture Imagery Program (NAIP) for the US\(^2\), and from the BD-ORTHO® 2.0 and ORTHO-HR® 1.0 databases from the IGN for France\(^3\). Land cover data originates from the National Land Cover Database (NLCD)\(^4\) for the US and from CESBIO\(^4\) for France. All elevation data comes from the NASA Shuttle Radar Topography Mission (SRTM)\(^5\). In addition, the dataset also includes traditional coarser resolution covariates: 19 bio-climatic rasters (30arcsec\(^2\)/pixel, i.e., 1km\(^2\)/pixel, from WorldClim [8]) and 8 pedologic rasters (250m\(^2\)/pixel, from SoilGrids [9]). The details of these rasters are given in Table 1.

Train-test split. The full set of occurrences was split in a training and testing set using a spatial block holdout procedure as illustrated in Figure 2. This limits the effect of spatial auto-correlation in the data [10]. Using this splitting procedure, a model cannot perform well by simply interpolating between training samples. The split was based on a global grid of 5km $\times$ 5km quadrats. 2.5% of these quadrats were randomly sampled and the observations falling in those formed the test set. 10% of those observations were used for the public leaderboard on Kaggle while the remaining 90% allowed to compute the private leaderboard providing the final results of the challenge. Similarly, another 2.5% of the quadrats were randomly sampled to provide an official validation set. The remaining quadrats and their associated observations were assigned to the training set.

Evaluation metric. For each occurrence in the test set, the goal of the task was to return a candidate set of species likely to be present at that location. Due to the presence-only [11] nature of the observation data used during the evaluation of the methods, for each location in the test set, we only have the knowledge of the presence of one species – the one observed – among the

\(^{2}\)https://www.fsa.usda.gov
\(^{3}\)https://geoservices.ign.fr
\(^{4}\)http://osr-cesbio.ups-tlse.fr/~oso/posts/2017-03-30-carte-s2-2016/
\(^{5}\)https://lpdaac.usgs.gov/products/srtmgl1v003/
Table 1
Summary of the low-resolution environmental variable rasters provided. The first 19 rows correspond to the bio-climatic variables from WorldClim [8]. The last 8 rows correspond to the pedologic variables from SoilGrid [9].

<table>
<thead>
<tr>
<th>Name</th>
<th>Description</th>
<th>Resolution</th>
</tr>
</thead>
<tbody>
<tr>
<td>bio_1</td>
<td>Annual Mean Temperature</td>
<td>30 arcsec</td>
</tr>
<tr>
<td>bio_2</td>
<td>Mean Diurnal Range (Mean of monthly (max temp - min temp))</td>
<td>30 arcsec</td>
</tr>
<tr>
<td>bio_3</td>
<td>Isothermality (bio_2/bio_7) (*100)</td>
<td>30 arcsec</td>
</tr>
<tr>
<td>bio_4</td>
<td>Temperature Seasonality (standard deviation *100)</td>
<td>30 arcsec</td>
</tr>
<tr>
<td>bio_5</td>
<td>Max Temperature of Warmest Month</td>
<td>30 arcsec</td>
</tr>
<tr>
<td>bio_6</td>
<td>Min Temperature of Coldest Month</td>
<td>30 arcsec</td>
</tr>
<tr>
<td>bio_7</td>
<td>Temperature Annual Range (bio_5-bio_6)</td>
<td>30 arcsec</td>
</tr>
<tr>
<td>bio_8</td>
<td>Mean Temperature of Wettest Quarter</td>
<td>30 arcsec</td>
</tr>
<tr>
<td>bio_9</td>
<td>Mean Temperature of Driest Quarter</td>
<td>30 arcsec</td>
</tr>
<tr>
<td>bio_10</td>
<td>Mean Temperature of Warmest Quarter</td>
<td>30 arcsec</td>
</tr>
<tr>
<td>bio_11</td>
<td>Mean Temperature of Coldest Quarter</td>
<td>30 arcsec</td>
</tr>
<tr>
<td>bio_12</td>
<td>Annual Precipitation</td>
<td>30 arcsec</td>
</tr>
<tr>
<td>bio_13</td>
<td>Precipitation of Wettest Month</td>
<td>30 arcsec</td>
</tr>
<tr>
<td>bio_14</td>
<td>Precipitation of Driest Month</td>
<td>30 arcsec</td>
</tr>
<tr>
<td>bio_15</td>
<td>Precipitation Seasonality (Coefficient of Variation)</td>
<td>30 arcsec</td>
</tr>
<tr>
<td>bio_16</td>
<td>Precipitation of Wettest Quarter</td>
<td>30 arcsec</td>
</tr>
<tr>
<td>bio_17</td>
<td>Precipitation of Driest Quarter</td>
<td>30 arcsec</td>
</tr>
<tr>
<td>bio_18</td>
<td>Precipitation of Warmest Quarter</td>
<td>30 arcsec</td>
</tr>
<tr>
<td>bio_19</td>
<td>Precipitation of Coldest Quarter</td>
<td>30 arcsec</td>
</tr>
<tr>
<td>orcdrc</td>
<td>Soil organic carbon content (g/kg at 15cm depth)</td>
<td>250 m</td>
</tr>
<tr>
<td>phihox</td>
<td>Ph x 10 in H2O (at 15cm depth)</td>
<td>250 m</td>
</tr>
<tr>
<td>ccesol</td>
<td>Cation exchange capacity of soil in cmolc/kg 15cm depth</td>
<td>250 m</td>
</tr>
<tr>
<td>bdticm</td>
<td>Absolute depth to bedrock in cm</td>
<td>250 m</td>
</tr>
<tr>
<td>clyppt</td>
<td>Clay (0-2 micro meter) mass fraction at 15cm depth</td>
<td>250 m</td>
</tr>
<tr>
<td>sltppt</td>
<td>Silt mass fraction at 15cm depth</td>
<td>250 m</td>
</tr>
<tr>
<td>sndppt</td>
<td>Sand mass fraction at 15cm depth</td>
<td>250 m</td>
</tr>
<tr>
<td>bldfie</td>
<td>Bulk density in kg/m3 at 15cm depth</td>
<td>250 m</td>
</tr>
</tbody>
</table>

Note that this evaluation metric does not try to correct the sampling bias [13] inherent to present-only observation data (linked to the density of population, etc.). The absolute value of the resulting figures should thus be taken with care. Nevertheless, this metric does allow to compare the different approaches and to determine which type of input data and of models are different ones which can actually be found all together at that point. To measure the precision of the predicted sets while accommodating with this limited knowledge, a simple set-valued classification [12] metric was chosen as the main evaluation criterion: top-30 error rate. Each observation \( i \) is associated with a single ground-truth label \( y_i \) corresponding to the observed species. For each observation, the submissions provided 30 candidate labels \( \hat{y}_{i,1}, \hat{y}_{i,2}, \ldots, \hat{y}_{i,30} \). The top-30 error rate is then computed using

\[
\text{Top-30 error rate} = \frac{1}{N} \sum_{i=1}^{N} e_i \quad \text{where} \quad e_i = \begin{cases} 
1 & \text{if } \forall k \in \{1, \ldots, 30\}, \hat{y}_{i,k} \neq y_i \\
0 & \text{otherwise}
\end{cases}
\]

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0 & \text{otherwise}
\end{cases}
\]
Course of the challenge. The training and test data were publicly shared in early March 2021 through the Kaggle platform⁶. Any research team wishing to participate in the evaluation could register on the platform and download the data. Each team could submit up to 3 submissions per day to compete on the public leaderboard. A submission (also called a run in the next sections) takes the form of a CSV file containing the top-30 predictions of the method being evaluated for all observations in the test set. For each submission, the top-30 error rate was first computed only on a subset of the test set to produce the public leaderboard which was visible to all the participants while the competition was still running. Once the submission phase was closed (mid-May), only 5 submissions per team were retained to compute the private leaderboard using the rest of the test set. These submissions were either hand-picked by the team or automatically chosen as the 5 best performing submissions on the public leaderboard. The participants could then see the final scores of all the other participants on the private leaderboard as well as their final ranking. Each participant was asked to provide a working note, i.e. a detailed report containing all technical information required to reproduce the results of the submissions. All LifeCLEF working notes were reviewed by at least two members of the LifeCLEF organizing committee to ensure a sufficient level of quality and reproducibility.

3. Baseline methods

Based on last year’s competition, three baselines were provided by the organizers of the challenge to serve as comparison references for the participants while developing their own methods. They consisted in:

- Top-30 most present species: a constant predictor returning always the same list of

⁶https://www.kaggle.com/c/geolifeclef-2021/
the most present species, i.e. the ones having the most occurrences in the training set.

- **RF on environmental variables**: a Random Forest model trained on environmental feature vectors only, i.e. on the 27 climatic and soil variables extracted at the position of the observation.

- **CNN on 6-channels patches**: this method [14] is the one that obtained the best result during GeoLifeCLEF 2020 competition [5]. It is based on a Convolution Neural Network trained on all the high-resolution image covariates, i.e. on 6-channel tensors composed of RGB-IR images, land cover image, and altitude image.

4. Participants and methods

Seven teams participated to the GeoLifeCLEF 2021 challenge and submitted a total of 26 submissions: University of Melbourne, DUTH (Democritus University of Thrace), CONABIO (Comisión Nacional para el Conocimiento y Uso de la Biodiversidad), UTFPR (Federal University of Technology – Paraná) as well as three participants for which we could not identify the affiliation and which we denote here, respectively, as Team Alpha, Team Beta and Team Gamma. The results are shown in Figure 3. Only the winning team (University of Melbourne) submitted a working notes paper with a detailed description of the used methodology and models trained [15]. We summarize hereafter the two main models they employed:

- **CNN on RGB patches (University of Melbourne - Run 1)**: this method is based on a Convolution Neural Network (ResNet50) trained on the RGB remote sensing images only. The model was first trained in an unsupervised way using MOCO on the dataset, a contrastive representation learning framework (for 20 epochs). The resulting model was then supervisely fine-tuned entirely (i.e., end-to-end) using 7 epochs of stochastic gradient descent (SGD) on the whole training set (validation set included). Three types of data augmentation were used to reduce overfitting: (i) random horizontal flip, (ii) random vertical flip, and (iii) RandAugment [16], an automated data augmentation optimizer.

- **Bi-modal CNN on RGB patches & altitude (University of Melbourne - Run 2)**: The CNN on RGB patches was combined with another CNN on the altitude patches (using the same ResNet50 architecture). This other CNN was also pre-trained in an unsupervised way similarly using MOCO as for the RGB-based model. The two models are then combined by concatenating the final bottleneck layer of the ResNet50 (i.e. the new layer contains 4096 neurons instead of 2048 in the mono-modal CNN on RGB patches) and the global model was fine-tuned as before on fewer epochs. Most data augmentation was removed during training.

5. Global competition results analysis

The global results of the GeoLifeCLEF 2021 are shown in Figure 3. Generally speaking, CNN-based models trained on high-resolution patches used in runs by University of Melbourne and Team Alpha as well as in the baseline CNN (high resolution patches) are very competitive and efficient compared to the traditional model (RF (environmental vectors)). This observation tends
to show that (i) important information explaining the species composition is contained in the high-resolution patches, and, (ii) convolutional neural networks are able to capture and exploit this information.

One question raised by the challenge is how to properly aggregate the different variables provided as input. Adding altitude data to the model (University of Melbourne - Run 2) provides an improvement in prediction accuracy backing the intuition that this variable is informative of the species distribution. However, aggregating all the variables does not mechanically lead to higher performance: CNN (high resolution patches) makes use of the additional land cover data but its performance is not as good as the two runs from University of Melbourne. It seems that it is important not to aggregate the features representation of those variables too early in the architectures of the networks: concatenation of higher-level features (University of Melbourne - Run 2) is more efficient than early aggregation (CNN (high resolution patches)).

### 6. Complementary analysis

In this section, we provide complementary analyses of the submitted results. These analyses are based on the complete predicted scores (the logits for every class of every test occurrence) gently provided by the University of Melbourne team [15] after the competition for both of their models. As the dataset used this year is the same as for GeoLifeCLEF 2020, some of the conclusions of last year’s overview paper [5] still hold and we refer the reader to it. Here, we focus on other aspects of the dataset not considered previously.
Comparison of set-valued classification metrics. As stated previously, the task considered here fits in the framework of set-valued classification [12]. A set of species is present at each location, the model tries to predict this set, however, we are only given a single of those present species to evaluate the performance of the model. This year, the top-30 error rate was chosen as the main evaluation metric. This metric however considers that, at each location, the same number of species will be present, implicitly assuming that the species richness is uniform over the geographical extent considered. To surpass this limitation, we can relax this constraint of predicting always sets of size $K$ by allowing to predict adaptive sets which size in average is equal to $K$. This type of set-valued classification is known as average-$K$ classification [17]. In Figure 4, we compare top-$K$ error rate to average-$K$ error rate for all values of $K$. Average-$K$ error rate is computed by finding the appropriate threshold on the scores which will result in sets of average size $K$ on the test set. We refer the reader to [17, 12] for more details. On the comparison plot, average-$K$ error rate can only improve upon top-$K$ error rate. However, in our case, both curves are indistinguishable. Meaning that either, average-$K$ does not capture the proper heterogeneity of the samples between each location [17] or that the predicted scores are not good enough to estimate it properly. We tried additional probability calibration procedures such as temperature scaling [18] to improve the average-$K$ predictions but it did not provide a significant edge. Further investigation is required to understand why average-$K$ set classifiers do not perform better on this dataset.

Kingdoms analysis. Most species are animals (16,328) and plants (13,035). Due to an omission during the construction of the dataset, other kingdoms are also represented in the US, accounting for 2,072 species (mainly Fungi) but representing only 1.5% of the observations. They are thus negligible. In Figure 5, we detail the performance of the bi-modal CNN of [15] for these different kingdoms. The main evaluation metric consists of all the kingdoms mixed together. For the separate kingdoms metrics, only the test points which observed species belonging to this specific kingdom are considered. If we directly compute the metrics on the top-30 sets
predicted as in Figure 5a, it would seem that the task is harder for animals than for plants. However, this gives a biased image of the difficulty of the task. A better approach is to compute separate top-30 sets for each kingdom by retaining only the species belonging to it. In this case, the predicted top-30 sets at the same location are thus different for each kingdom. The resulting error rates are shown in Figure 5b. In that case, animals, although containing more species and fewer observations, seem easier to predict than plants. Surprisingly, although the dataset contains few (28,573) observations from other kingdoms, they seem to be easier to predict. This is actually likely to be simply an artefact of the metric: as they are relatively few species compared to the other kingdoms, taking the top-30 most probable species is probably a too large and not very informative set. This analysis suggests that predicting separate top-\(K\) sets for each kingdom and use these sets to derive a new evaluation metric is worth considering. Note that, this might not be as direct as it seems. Indeed, it could be important to use different values of \(K\) for the different kingdoms depending of the number of species they contain and the likelihood to find several of them present at the same location. This would thus require some tweaking.

**Late-fusion of models.** We test the complementarity of the predictions made by the different methods by aggregating their output (the logits) in a late-fusion manner: their outputs are averaged before computing the top-30 most confident species. Here, we focus on the runs submitted by team University of Melbourne as well as the baseline CNN (high resolution patches). The results are shown in Figure 6. As explained previously, one important difference between the baseline and the runs submitted is the way the variables are aggregated. Surprisingly, fusing the predictions of the baseline method with either the uni-modal or bi-modal model from University of Melbourne actually results in top-30 sets of significantly similar accuracy. This seems to suggest that, beyond the aggregation method, the unsupervised pre-training used by University of Melbourne might be particularly helpful for RGB data but less for altitude data.
which is already fairly well exploited by the baseline. Another interesting observation is that combining the uni-modal and bi-modal models from [15] provides a little additional gain of about 2% This might come up as a surprise: the bi-modal model consists essentially of the uni-modal model with additional altitude data. This additional gain might come from three factors:

- the slight differences in the training of uni-modal and bi-modal models, i.e. data augmentation and training time;
- the late aggregation of the altitude data with RGB images might not be perfectly suited and adding a second layer of aggregation via late-fusion might allow to recover it additional complementary information;
- the variance reduction due to ensembling of similar models with high bias.

Finally, late-fusing all three models provides yet another non-negligible gain, highlighting the complementary of those different models. Further investigation could understand better where this complementary originates and help build better models.

7. Conclusion

The main challenges raised by this year’s challenge are two folds:

- How to properly aggregate the different covariates provided?
- How much complementary or redundant is the information contained in the high-resolution patches to the one captured from the bioclimatic and soil variables?
Moreover, there remains considerable room for improvement on this challenge as the winning solution does not make use of all the different patches provided and its top-30 error rate is still high, near 75% error rate.

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