# Guidelines for the validation of machine learning predictions of species interactions

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- 1. The prediction of species interactions is gaining momentum as a way to circumvent limitations in data volume. Yet, ecological networks are challenging to predict because they are typically small and sparse. Dealing with extreme class imbalance is a challenge for most binary classifiers, and there are currently no guidelines as to how predictive models can be trained for this specific problem.
- 2. Using simple mathematical arguments and numerical experiments in which a variety of classifiers (for supervised learning) are trained on simulated networks, we develop a series of guidelines related to the choice of measures to use for model selection, and the degree of unbiasing to apply to the training dataset.
- 3. Neither classifier accuracy nor the ROC-AUC are informative measures for the performance of interaction prediction. PR-AUC is a fairer assessment of performance. In some cases, even standard measures can lead to selecting a more biased classifier because the effect of connectance is strong. The amount of correction to apply to the training dataset depends on network connectance, on the measure to be optimized, and only weakly on the classifier.
- 4. These results reveal that training machines to predict networks is a challenging task, and that in virtually all cases, the composition of the training set needs to be experimented on before performing the actual training. We discuss these consequences in the context of the low volume of data.

- Ecological networks are a backbone for key ecological and evolutionary processes; yet enumerating all of
- the interactions they contain is a daunting task, as it scales with  $S^2$ , i.e. the squared species richness
- TODO. Recent contributions to the field of ecological network prediction (Pichler et al., 2020; Strydom et
- 4 al., 2021; Becker2021OptPre?) highlight that although interactions can be predicted by adding
- ecologically relevant information (in the form of, e.g. traits), we do not have robust guidelines as to how
- 6 the predictive ability of these models should be evaluated, nor about how the models should be trained.
- 7 Here, by relying on simple derivations and a series of simulations, we formulate a number of such
- 8 guidelines, specifically for the case of binary classifiers derived from thresholded values. Specifically, we
- 9 conduct an investigation of the models in terms of their skill (ability to make the right prediction), bias
- 10 (trends towards systematically over-predicting one class), class imbalance (the relative number of cases
- representing interactions), and show how these effects interact. We conclude on the fact that models with
- the best interaction-scale predictive score do not necessarily result in the most accurate representation of
- 13 the network.
- 14 The prediction of ecological interactions shares conceptual and methodological issues with two fields in
- biology: species distribution models (SDMs), and genomics. SDMs suffers from issues affecting
- interactions prediction, namely low prevalence (due to sparsity of observations/interactions) and data
- aggregation (due to bias in sampling some locations/species). In previous work, Allouche et al. (2006)
- suggested that  $\kappa$  was a better test of model performance than the True Skill Statistic (TSS; which we refer
- to as Youden's informedness); these conclusions were later criticized by Somodi et al. (2017), who
- emphasized that informedness' is affected both by prevalence and bias. Although this work offers
- 21 recommendations about the comparison of models, it doesn't establishes baselines or good practices for
- training on imbalanced ecological data, or ways to remedy the imbalance. Steen et al. (2021) show that,
- when applying spatial thinning (a process that has no analogue in networks), the best approach to train
- ML-based SDMs varies according to the balancing of the dataset, and the evaluation measures used. This
- suggests that there is no single "recipe" that is guaranteed to give the best model. By contrast to networks,
- SDMs have the advantage of being able to both thin datasets to remove some of the sampling bias (e.g.
- 27 Inman et al., 2021), but also to create pseudo-absences to inflate the number of supposed negatives in the
- dataset (e.g. Iturbide et al., 2015).
- 29 An immense body of research on machine learning application to life sciences is focused on genomics
- 30 (which has very specific challenges, see a recent discussion by Whalen et al., 2021); this sub-field has

- 31 generated recommendations that do not necessarily match the current best-practices for SDMs, and
- therefore hint at the importance of domain-specific guidelines. Chicco & Jurman (2020) suggest using
- Matthews correlation coefficient (MCC) over  $F_1$ , as a protection against over-inflation of predicted results;
- Delgado & Tibau (2019) advocate against the use of Cohen's  $\kappa$ , again in favor of MCC, as the relative
- nature of  $\kappa$  means that a worse classifier can be picked over a better one; similarly, Boughorbel et al.
- 36 (2017) recommend MCC over other measures of performance for imbalanced data, as it has more
- desirable statistical properties. More recently, Chicco et al. (2021) temper the apparent supremacy of the
- $^{38}$  MCC, by suggesting it should be replaced by Youden's informedness (also known as J, bookmaker's
- 39 accuracy, and the True-Skill Statistic) when the imbalance in the dataset may not be representative of the
- 40 actual imbalance.
- Species interaction networks are often under-sampled (Jordano, 2016b, 2016a), and this under-sampling is
- 42 structured taxonomically (Beauchesne et al., 2016), structurally (de Aguiar et al., 2019) and spatially
- (Poisot, Bergeron, et al., 2021; Wood et al., 2015). As a consequence, networks suffer from data
- deficiencies both within and between datasets. This implies that the comparison of classifiers across
- space, when undersampling varies locally (see e.g. McLeod et al., 2021) is non-trivial. Furthermore, the
- 46 baseline value of classifiers performance measures under various conditions of skill, bias, and prevalence,
- 47 has to be identified to allow researchers to evaluate whether their interaction prediction model is indeed
- 48 learning. Taken together, these considerations highlight three specific issues for ecological networks.
- 49 First, what values of performance measures are indicative of a classifier with no skill? This is particularly
- 50 important as it can evaluate whether low prevalence can lull us into a false sense of predictive accuracy.
- 51 Second, independently of the question of model evaluation, is low prevalence an issue for training or
- testing, and can we remedy it? Finally, because the low amount of data on interaction makes a lot of
- imbalance correction methods (see e.g. Branco et al., 2015) hard to apply, which indicators can be
- optimized by sacrificing least amount of positive interaction data?
- 55 It may sound counter-intuitive to care so deeply about how good a classifier with no-skill is, as by
- definition, is has no skill. The necessity of this exercise has its roots in the paradox of accuracy: when the
- desired class ("two species interact") is rare, a model that gets less ecologically performant by only
- predicting the opposite class ("these two species do not interact") sees its accuracy increase; because most
- of the guesses have "these two species do not interact" as a correct answer, a model that never predicts
- 60 interactions would be right an overwhelming majority of the time; it would also be utterly useless. Herein

- 61 lies the core challenge of predicting species interactions: the extreme imbalance between classes makes
- the training of predictive models difficult, and their validation even more so as we do not reliably know
- which negatives are true. The connectance (the proportion of realized interactions, usually the number of
- 64 interactions divided by the number of species pairs) of empirical networks is usually well under 20%, with
- larger networks having a lower connectance (MacDonald et al., 2020), and therefore being increasingly
- 66 difficult to predict.

# A primer on binary classifier evaluation

- <sup>68</sup> Binary classifiers, which it to say, machine learning algorithms whose answer is a categorical value, are
- 69 usually assessed by measuring properties of their confusion matrix, i.e. the contingency table reporting
- 70 true/false positive/negative hits. A confusion matrix is laid out as

$$\begin{pmatrix} tp & fp \\ fn & tn \end{pmatrix}$$
.

- In this matrix, tp is the number of times the model predicts an interaction that exists in the network (true
- positive), fp is the number of times the model predicts an interaction that does not exist in the network
- <sub>73</sub> (false positive), fn is the number of times the model fails to predict an interaction that actually exists in the
- network (false negatives), and to is the number of times the model correctly predicts that an interaction
- does not exist (true negatives). From these values, we can derive a number of measures of model
- performance (see Strydom et al., 2021 for a review of their interpretation in the context of networks). At a
- coarse scale, a classifier is accurate when the trace of the matrix divided by the sum of the matrix is close
- to 1, with other measures informing us on how the predictions fail.
- 79 There is an immense diversity of measures to evaluate the performance of classification tasks (Ferri et al.,
- 2009). Here we will focus on five of them with high relevance for imbalanced learning (He & Ma, 2013).
- The choice of metrics with relevance to class-imbalanced problems is fundamental, because as Japkowicz
- 82 (2013) unambiguously concluded, "relatively robust procedures used for unskewed data can break down
- miserably when the data is skewed." Following Japkowicz (2013), we focus on two ranking metrics (the
- areas under the Receiver Operating Characteristic and Precision Recall curves), and three threshold

- metrics ( $\kappa$ , informedness, and MCC; we will briefly discuss  $F_1$  but show early on that it has undesirable
- 86 properties).
- The  $\kappa$  measure (Landis & Koch, 1977) establishes the extent to which two observers (the network and the
- prediction) agree, and is measured as

$$2\frac{tp\times tn-fn\times fp}{(tp+fp)\times (fp+tn)+(tn+fp)\times (tn+fn)}\,.$$

- Informedness (Youden, 1950) (also known as bookmaker informedness or the True Skill Statistic) is
- TPR + TNR 1, where TPR = tp/(tp + fn) and TNR = tn/(tn + fp). Informedness can be used to find
- 91 the optimal cutpoint in thresholding analyses (Schisterman et al., 2005); indeed, the maximal
- 92 informedness corresponds to the point on the ROC curve that is closest to the perfect classifier point. The
- 93 formula for informedness is

$$\frac{tp}{tp+fn} + \frac{tn}{tn+fp} - 1.$$

94 The MCC is defined as

$$\frac{tp \times tn - fn \times fp}{\sqrt{(tp + fp) \times (tp + fn) \times (tn + fp) \times (tn + fn)}}.$$

- Finally,  $F_1$  is the harmonic mean of precision (the chance that interaction was correctly detected as such)
- and sensitivity (the ability to correctly classify interactions), and is defined as

$$2\frac{tp}{2\times tp + fp + fn}.$$

- 97 A lot of binary classifiers are built by using a regressor (whose task is to guess the value of the interaction,
- 98 and can therefore return a value considered to be a pseudo-probability); in this case, the optimal value
- below which predictions are assumed to be negative (i.e. the interaction does not exist) can be determined
- by picking a threshold maximizing some value on the ROC or the PR curve. The area under these curves
- (ROC-AUC and PR-AUC henceforth) give ideas on the overall goodness of the classifier, and the ideal
- threshold is the point on these curves that minimizes the tradeoff represented in these curves. Saito &

Rehmsmeier (2015) established that the ROC-AUC is biased towards over-estimating performance for imbalanced data; on the contrary, the PR-AUC is able to identify classifiers that are less able to detect 104 positive interactions correctly, with the additional advantage of having a baseline value equal to 105 prevalence. Therefore, it is important to assess whether these two measures return different results when applied to ecological network prediction. The ROC curve is defined by the false positive rate on the x axis, 107 and the true positive rate on the y axis, and the PR curve is defined by the true positive rate on the x axis, 108 and the positive predictive value on the y axis. By comparison with the previous paragraph, it is obvious 109 that  $F_1$  and MCC have ties to the PR curve (being close to the expected PR-AUC), and that informedness 110 has ties to the ROC curve (whereby the threshold maximizing informedness is also the point of maximal 111 inflection on the ROC curve). One important difference between ROC and PR is that the later does not 112 prominently account for the size of the true negative compartments: in short, it is more sensitive to the correct positive predictions. In a context of strong imbalance, PR-AUC is therefore a more stringent test of 114 model performance. 115

## Baseline values for the threshold metrics

In this section, we will assume a network of connectance  $\rho$ , *i.e.* having  $\rho S^2$  interactions (where S is the species richness), and  $(1-\rho)S^2$  non-interactions. Therefore, the vector describing the *true* state of the network (assumed to be an unweighted, directed network) is a column vector  $\mathbf{o}^T = [\rho, (1-\rho)]$  (we can safely drop the  $S^2$  terms, as we will work on the confusion matrix, which ends up expressing *relative* values). We will apply skill and bias to this matrix, and measure how a selection of performance metrics respond to changes in these values, in order to assess their suitability for model evaluation.

#### Confusion matrix with skill and bias

In order to write the values of the confusion matrix for a hypothetical classifier, we need to define two characteristics: its skill, and its bias. Skill, here, refers to the propensity of the classifier to get the correct answer (*i.e.* to assign interactions where they are, and to not assign them where they are not). A no-skill classifier guesses at random, *i.e.* it will guess interactions with a probability  $\rho$ . The predictions of a no-skill classifier can be expressed as a row vector  $\mathbf{p} = [\rho(1-\rho)]$ . The confusion matrix  $\mathbf{M}$  for a no-skill classifier is given by the element-wise (Hadamard, outer) product of these vectors  $\mathbf{o} \odot \mathbf{p}$ , *i.e.* 

$$\mathbf{M} = \begin{pmatrix} \rho^2 & \rho(1-\rho) \\ (1-\rho)\rho & (1-\rho)^2 \end{pmatrix}.$$

In order to regulate the skill of this classifier, we can define a skill matrix **S** with diagonal elements equal to s, and off-diagonal elements equal to s, and re-express the skill-adjusted confusion matrix as  $\mathbf{M} \odot \mathbf{S}$ , *i.e*.

$$\begin{pmatrix} \rho^2 & \rho(1-\rho) \\ (1-\rho)\rho & (1-\rho)^2 \end{pmatrix} \odot \begin{pmatrix} s & (1-s) \\ (1-s) & s \end{pmatrix}.$$

When s = 0,  $Tr(\mathbf{M}) = 0$  (the classifier is *always* wrong), when s = 0.5, the classifier is no-skill and guesses at random, and when s = 1, the classifier is perfect.

The second element we can adjust in this hypothetical classifier is its bias, specifically its tendency to over-predict interactions. Like above, we can do so by defining a bias matrix  $\mathbf{B}$ , where interactions are over-predicted with probability b, and express the final classifier confusion matrix as  $\mathbf{M} \odot \mathbf{S} \odot \mathbf{B}$ , *i.e.* 

$$\begin{pmatrix} \rho^2 & \rho(1-\rho) \\ (1-\rho)\rho & (1-\rho)^2 \end{pmatrix} \odot \begin{pmatrix} s & (1-s) \\ (1-s) & s \end{pmatrix} \odot \begin{pmatrix} b & b \\ (1-b) & (1-b) \end{pmatrix}.$$

The final expression for the confusion matrix in which we can regulate the skill and the bias is

$$\mathbf{C} = \begin{pmatrix} s \times b \times \rho^2 & (1-s) \times b \times \rho(1-\rho) \\ (1-s) \times (1-b) \times (1-\rho)\rho & s \times (1-b) \times (1-\rho)^2 \end{pmatrix}.$$

In all further simulations, the confusion matrix **C** is transformed so that it sums to unity, *i.e.* the entries are the *proportions* of guesses.

#### What are the baseline values of performance measures?

In this section, we will change the values of b, s, and  $\rho$ , and report how the main measures discussed in the introduction (MCC,  $F_1$ ,  $\kappa$ , and informedness) respond. Before we do so, it is important to explain why

the sum of the confusion matrix. For a no-skill, no-bias classifier, accuracy is equal to  $\rho^2 + (1 - \rho)^2$ ; for 145  $\rho = 0.05$ , this is  $\approx 0.90$ , and for  $\rho = 0.01$ , this is equal to  $\approx 0.98$ . In other words, the values of accuracy are 146 high enough to be uninformative (for  $\rho$  small,  $\rho^2 \ll (1-\rho)^2$ ). More concerning is the fact that introducing bias changes the response of accuracy in unexpected ways. Assuming a no-skill classifier, the numerator 148 of accuracy becomes  $b\rho^2 + (1-b)(1-\rho)^2$ , which increases when b is low, which specifically means that at 149 equal skill, a classifier that under-predicts interactions will have higher accuracy than an un-biased 150 classifier (because the value of accuracy is dominated by the size of tn, which will increase). These issues 151 are absent from balanced accuracy, but should nevertheless lead us to not report accuracy as the primary 152 measure of network prediction success; moving forward, we will focus on other measures. 153 In order to examine how MCC,  $F_1$ ,  $\kappa$ , and informedness change w.r.t. the imbalance, skill, and bias, we 154 performed a grid exploration of the values of logit(s) and logit(b) linearly from -10 to 10; logit(x) = -10155 means that x is essentially 0, and logit(x) = 10 means it is essentially 1 – this choice was motivated by the fact that most responses are non-linear with regards to bias and skill. The values or  $\rho$  were taken linearly 157 in [0, 0.5], which is within the range of connectance for species interaction networks. Note that at this 158 point, there is no network model to speak of; the confusion matrix we discuss can be obtained for any 159 classification task. Based on the previous discussion, the desirable properties for a measure of classifier 160 success should be: an increase with classifier skill, especially at low bias; a hump-shaped response to bias, 161 especially at high skill, and ideally centered around logit(b) = 0; an increase with prevalence up until 162 equiprevalence is reached.

we will not focus on accuracy too much. Accuracy is the number of correct predictions ( $Tr(\mathbf{C})$ ) divided by

#### [Figure 1 about here.]

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In fig. 1, we show that none of the four measures satisfy all the considerations at once:  $F_1$  increases with skill, and increases monotonously with bias; this is because  $F_1$  does not account for true negatives, and the increase in positive detection masks the over-prediction of interactions. Informedness varies with skill, reaching 0 for a no-skill classifier, but is entirely unsensitive to bias. Both MCC and  $\kappa$  have the same behavior, whereby they increase with skill.  $\kappa$  peaks at increasing values of biass for increasing skill, *i.e.* is likely to lead to the selection of a classifier that over-predicts interactions. By contract, MCC peaks at the same value, regardless of skill, but this value is not logit(b) = 0: unless at very high classifier skill, MCC risks leading to a model that over-predicts interactions. In fig. 2, we show that all measures except  $F_1$  give

a value of 0 for a no-skill classifier, and are forced towars their correct maximal value when skill changes
(i.e. a more connected networks will have higher values for a skilled classifierd, and lower values for a
classifier making mostly mistakes).

## [Figure 2 about here.]

These two analyses point to the following recommendations: MCC is indeed more appropriate than  $\kappa$ , as 177 although sensitive to bias, it is sensitive in a consistent way. Informedness is appropriate at discriminating 178 between different skills, but confounded by bias. As both of these measures bring valuable information on 179 the model behavior, we will retain them for future analyses.  $F_1$  is increasing with bias, and should not be 180 prioritized to evalue the performance of the model. The discussion of sensitivity to bias should come with 181 a domain-specific caveat: although it is likely that interactions documented in ecological networks are 182 correct, a lot of non-interactions are simply unobserved; as predictive models are used for data-inflation 183 (i.e. the prediction of new interactions), it is not necessarily a bad thing in practice to select models that predict more interactions than the original dataset, because the original dataset misses some interactions. 185 Furthermore, the weight of positive interactions could be adjusted if some information about the extent of 186 undersampling exists (e.g. Branco et al., 2015). In a recent large-scale imputation of interactions in the 187 mammal-virus networks, Poisot, Ouellet, et al. (2021) for example estimated that 93% of interactions are 188 yet to be documented. 189

# Numerical experiments on training strategy

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In the following section, we will generate random bipartite networks (this works without loss of generality 191 on unipartite networks), and train four binary classifiers (as well as an ensemble model using the sum of 192 ranged outputs from the component models) on 30% of the interaction data. In practice, testing usually 193 uses 70% of the total data; for ecological networks, where interactions are sparse and the number of 194 species is low, this may not be the best solution, as the testing set becomes constrained not by the 195 proportion of interactions, but by their number. Preliminary experiments using different splits revealed no 196 qualitative change in the results. Networks are generated by picking a random infectiousness trait  $v_i$  for 197 100 species (from a beta distribution  $B(\alpha = 6, \beta = 8)$  distribution), and a resistance trait  $h_i$  for 100 species 198 (from  $B(\alpha = 2, \beta = 8)$  distribution). There is an interaction between i and j when

 $v_i - \xi/2 \le h_i \le v_i + \xi/2$ , where  $\xi$  is a constant regulating the connectance of the network (there is an almost 1:1 relationship between  $\xi$  and connectance), and varies uniformly in [0.05, 0.35]. This model gives 201 fully interval networks that are close analogues to the bacteria-phage model of Weitz et al. (2005), with 202 both a modular structure and a non-uniform degree distribution. This dataset is easy for almost any 203 algorithm to learn: when trained with features  $[v_i, h_j, abs(v_i, h_j)]^T$  to predict the interactions between i 204 and j, all four models presented below were able to reach almost perfect predictions all the time (data not 205 presented here) - this is in part because the rule (there is maximum value of the distance between traits 206 for which there is an interaction) is fixed for all interactions. In order to make the problem more difficult 207 to solve, we use  $[v_i, h_i]$  as a feature vector (i.e. the traits on which the models are trained), and therefore 208 the models will have to uncover that the rule for interaction is  $abs(v_i, h_i) \le \xi$ . The models therefore all 209 have the following form, where  $i_{i,j}$  is an interaction from species i to species j:

$$\begin{bmatrix} i_{1,1} \\ i_{1,2} \\ \vdots \\ i_{m,n-1} \\ i_{m,n} \end{bmatrix} \propto \begin{bmatrix} v_1 & h_1 \\ v_1 & h_2 \\ \vdots & \vdots \\ v_m & h_{n-1} \\ v_m & h_n \end{bmatrix}$$

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The training sample is composed of 30% of the  $10^4$  possible entries in the network, i.e. n = 3000. Out of these interactions, we pick a proportion  $\nu$  (the training set balance) to be positive, so that the training set 212 has  $\nu n$  interactions, and  $(1-\nu)n$  non-interactions. We vary  $\nu$  uniformly in ]0,1[. This allows to evaluate 213 how the measures of binary classification performance respond to artificially rebalanced dataset for a 214 given network connectance. The rest of the dataset (n = 7000 pairs of species) is used as a testing set, on 215 which all furher measures are calculated. Note that although the training set is balanced, the testing set is 216 not, and retains (part of) the imbalance of the original data. 217 The dataset used for numerical experiments is composed of 64000 such  $(\xi, \nu)$  pairs, on which four 218 machines are trained: a decision tree regressor, a boosted regression tree, a ridge regressor, and a random 219 forest regressor. All models were taken from the MLJ. jl package (Blaom et al., 2020; Blaom & Vollmer, 220 2020) in Julia 1.7 (Bezanson et al., 2017). All machines use the default parameterization; this is an obvious 221 deviation from best practices, as the hyperparameters of any machine require training before its 222 application on a real dataset. As we use 64000 such datasets, this would require 256000 unique instances 223

of tweaking the hyperparameters, which is not realistic. Therefore, we assume that the default parameterizations are comparable across networks. All machines return a quantitative prediction, usually 225 (but not necessarily) in [0,1], which is proportional (but not necessarily linearly) to the probability of an 226 interaction between i and j. Nevertheless, the ROC-AUC and PR-AUC (and therefore the thresholds) can be measured by integrating over the domain of the values return by each machine. 228 In order to pick the best adjacency matrix for a given trained machine, we performed a thresholding 229 approach using 500 steps on predictions from the testing set, and picking the threshold that maximized Youden's informedness. During the thresholding step, we measured the area under the receiver operating 231 characteristic (ROC-AUC) and precision-recall (PR-AUC) curves, as measures of overall performance over 232 the range of returned values. We report the ROC-AUC and PR-AUC, as well as a suite of other measures as introduced in the next section, for the best threshold. The ensemble model was generated by summing the 234 predictions of all component models on the testing set (ranged in [0,1]), then put through the same 235 thresholding process. The complete code to run the simulations is available at 10.17605/OSF. IO/JKEWD. After the simulations were completed, we removed all runs (i.e. pairs of  $\xi$  and  $\nu$ ) for which at least one of 237 the following conditions was met: the accuracy was 0, the true positive or true negative rates were 0, the 238 connectance was larger than 0.25. This removes both the obviously failed model runs, and the networks 239 that are more densely connected compared to the connectance of empirical food webs (and are therefore 240 less difficult to predict, being less imbalanced; preliminary analyses of data with a connectance larger than 241 0.3 revealed that all machines reached consistently high performance).

# Effect of training set balance on performance

In fig. 3, we present the response of two thresholding measures (PR-AUC and ROC-AUC) and two ranking
measures (Informedness and MCC) to a grid of 35 values of training set balance, and 35 values of
connectance, for the four component models as well as the ensemble. ROC-AUC is always high, and does
not vary with training set balance. On the other hand, PR-AUC shows very strong responses, increasing
with training set balance. It is notable here that two classifiers that seemed to be performing well (Decision
Tree and Random Forest) based on their MCC are not able to reach a high PR-AUC even at higher
connectances. All models reached a higher performance on more connected networks, and using more
balanced training sets. In all cases, informedness was extremely high, which is an expected consequence

of the fact that this is the value we optimized to determine the cutoff. MCC increased with training set
balance, although this increase became less steep with increasing connectance. Three of the models (kNN,
decision tree, and random forest) only increased their PR-AUC sharply when the training set was heavily
imbalanced towards more interactions. Interestingly, the ensemble almost always outclassed its
component models. For larger connectances (less difficult networks to predict, as they are more balanced),
MCC and informedness stared decreasing when the training set bias got too close to one, suggesting that a
training set balance of 0.5 may often be appropriate if these measures are the one to optimize.

## [Figure 3 about here.]

Based on the results presented in fig. 3, it seems that informedness and ROC-AUC are not necessarily able to discriminate between good and bad classifiers (although this result may be an artifact for informedness, as it has been optimized when thresholding). On the other hand, MCC and PR-AUC show a strong response to training set balance, and may therefore be more useful at model comparison.

# Required amount of positives to get the best performance

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The previous results revealed that the measure of classification performance responds both to the bias in 265 the training set and to the connectance of the network; from a practical point of view, assembling a 266 training set requires to withold positive information, which in ecological networks are very scarce (and 267 typically more valuable than negatives, on which there is a doubt). For this reason, across all values of 268 connectance, we measured the training set balance that maximized a series of performance measures. 269 When this value is high, the training set needs to skew more positive in order to get a performant model; 270 when this value is about 0.5, the training set needs to be artificially balanced to optimize the model 271 performance. These results are presented in fig. 4. 272

#### [Figure 4 about here.]

The more "optimistic" measures (ROC-AUC and informedness) required a biasing of the dataset from about 0.4 to 0.75 to be maximized, with the amount of bias required decreasing only slightly with the connectance of the original network. MCC and PR-AUC required values of training set balance from 0.75 to almost 1 to be optimized, which is in line with the results of the previous section, *i.e.* they are more

stringent tests of model performance. These results suggest that learning from a dataset with very low
connectance can be a different task than for more connected networks: it becomes increasingly important
to capture the mechanisms that make an interaction *exist*, and therefore having a slightly more biased
training dataset might be beneficial. As connectance increases, the need for biased training sets is less
prominent, as learning the rules for which interactions *do not* exist starts gaining importance.

## [Figure 5 about here.]

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When trained at their optimal training set balance, connectance still had a significant impact on the
performance of some machines fig. 5. Notably, Decision Tree, Random Forest, and Ridge Regression had
low values of PR-AUC. In all cases, the Boosted Regression Tree was reaching very good predictions
(especially for connectances larger than 0.1), and the ensemble was almost always scoring perfectly. This
suggests that all the models are biased in different ways, and that the averaging in the ensemble is able to
correct these biases. We do not expect this last result to have any generality, and provide a discussion of a
recent example in which the ensemble was performing worse than its components models.

# Do better classification accuracy result in more realistic networks?

In this last section, we generate a network using the same model as before, with  $S_1, S_2 = 50, 80$  species, a connectance of  $\approx 0.16$  ( $\xi = 0.19$ ), and a training set balance of 0.5, as fig. 4 suggests this is the optimal training set balance for this range of connectance. The prediction made on the complete dataset is presented in fig. 6. Visualizing the results this way highlights the importance of exploratory data analysis: whereas all models return a network with interactions laying mostly on the diagonal (as expected), the Ridge Regression is quite obviously biased. Despite this, we can see that the ensemble is close to the initial dataset.

# [Figure 6 about here.]

The trained models were then thresholded (again by optimising informedness), and their predictions transformed back into networks for analysis; specifically, we measured the connectance, nestedness (REF), modularity (REF), asymmetry, and network dissimilarity (REF). This process was repeated 250

times, and the results are presented in tbl. 1. The random forest model is an interesting instance here: it produces the network that looks the most like the original dataset, despite having a very low PR-AUC, 304 suggesting it hits high recall at the cost of low precision. Although the ensemble was able to reach a very 305 high PR-AUC (and a very high ROC-AUC), this did not necessarily translate into more accurate reconstructions of the structure of the network. This result bears elaborating. Measures of model 307 performance capture how much of the interactions and non-interactions are correctly identified. As long 308 as these predictions are not perfect, some interactions will be predicted at the "wrong" position in the 309 network; these measures cannot describe the structural effect of these mistakes. On the other hand, 310 measures of network structure can have the same value with interactions that fall at drastically different 311 positions; this is in part because a lot of these measures covary with connectance, and in part because as 312 long as these values are not 0 or their respective maximum, there is a large number of network configurations that can have the same value. That ROC-AUC is consistently larger than PR-AUC may be a 314 case of this measure masking models that are not, individually, strong predictors (Jeni et al., 2013). 315

Table 1: Values of four performance metrics, and five network structure metrics, for 500 independent predictions similar to the ones presented in fig. 6. The values in **bold** indicate the best value for each column (including ties). Because the values have been rounded, values of 1.0 for the ROC-AUC column indicate an average  $\geq 0.99$ .

Model	MCC	Inf.	ROC-AUC	PR-AUC	Conn.	η	Q
Decision tree	0.85	0.92	0.97	0.12	0.21	0.76	0.31
BRT	0.90	0.90	0.98	0.86	0.23	0.82	0.27
Random Forest	0.90	0.96	1.00	0.27	0.20	0.72	0.32
Ridge Regression	0.80	0.91	0.95	0.58	0.24	1.0	0.18
Ensemble	0.88	0.94	1.00	0.96	0.20	0.75	0.31
Data					0.18	0.66	0.34
<del></del>							

# Guidelines for the assesment of network predictive models

We establish that due to the low prevalence of interactions, even poor classifiers applied to food web data
will reach a high accuracy; this is because the measure is dominated by the accidentally correct
predictions of negatives. On simulated confusion matrices with ranges of imbalance that are credible for

ecological networks, MCC had the most desirable behavior, and informedness is a linear measure of classifier skill. By performing simulations with four models and an ensemble, we show that informedness 321 and ROC-AUC are consistently high on network data, and that MCC and PR-AUC are more accurate 322 measures of the effective performance of the classifier. Finally, by measuring the structure of predicted networks, we highlight an interesting paradox: the models with the best performance measures are not 324 the models with the closest reconstructed network structure. We discuss these results in the context of 325 establishing guidelines for the prediction of ecological interactions. 326 TODO informedness and accuracy should be easy to beat, make sure the model is better than them! The results presented here highlight an interesting paradox: although the Random Forest was ultimately 328 able to get a correct estimate of network structure tbl. 1, it ultimately remains a poor classifier, as 329 evidenced by its low PR-AUC. This suggests that the goal of predicting interactions and predicting networks may not be solvable in the same way – of course a perfect classifier of interactions would make a 331 perfect network prediction; but even the best scoring predictor of interactions (the ensemble model) had 332 not necessarily the best prediction of network structure. The tasks of predicting networks structure and of 333 predicting interactions within networks are essentially two different ones. For some applications (.e.g. 334 comparison of network structure across gradients), one may care more about a robust estimate of the 335 structure, at the cost at putting some interactions at the wrong place. For other applications (e.g. 336 identifying pairs of interacting species), one may conversely care more about getting as many pairs right, 337 even though the mistakes accumulate in the form of a slightly worse estimate of network structure. How 338 these two approaches can be reconciled is undoubtedly a task for further research. Despite this apparent 339 tension at the heart of the predictive exercise, we can use the results presented here to suggest a number of guidelines. 341 First, because we should have more trust in reported interactions than in reported absences of interactions, 342 we can draw on previous literature to recommend informedness as a measure to decide on a threshold (Chicco et al., 2021); this being said, because informedness is insensitive to bias, the model performance is 344 better evaluated through the use of MCC fig.  $\ref{MCC}$ . Because  $F_1$  is monotonously sensitive to classifier bias 345 fig. 1 and network connectance fig. 2, MCC should be prefered as a measure of model evaluation. 346 Second, because the PR-AUC responds more to network connectance fig. 5 and training set imbalance fig. ??, it should be used as a measure of model performance over the ROC-AUC. This is not to say that 348 ROC-AUC should be discarded (in fact, a low ROC-AUC is a sign of an issue with the model), but that its

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interpretation should be guided by the PR-AUC value. Specifically, a high ROC-AUC is not informative, as
    it can be associated to a low PR-AUC (see e.g. Random Forest in tbl. 1) This again echoes
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    recommendations from other fields (Jeni et al., 2013; Saito & Rehmsmeier, 2015).
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    Thirdly, regardless of network connectance, maximizing informedness required a training set balance of
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    about 0.5, and maximizing the MCC required a training set bias of 0.7 and more. This has an important
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    consequence in ecological networks, for which the pool of positive cases (interactions) to draw from is
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    typically small: the most parsimonious measure (i.e. the one requiring to discard the least amount of
    information to train the model) will give the best validation potential, and is probably the informedness
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    (maximizing informedness is the generally accepted default for imbalanced classification; Schisterman et
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    al., 2005).
    Finally, it is noteworthy that the ensemble model was systematically better than the component models;
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    even when the models were individually far form perfect, the ensemble was able to leverage the different
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    biases expressed by the models to make an overall more accurate prediction. We do not expect that
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    ensembles will always be better than single models. In a recent multi-model comparison,
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    (Becker2021OptPre?) found that the ensemble was not the best model. There is no general conclusion to
    draw from this besides reinforcing the need to be pragmatic about which models should be included in the
    ensemble, or whether to use an ensemble at all. In a sense, the surprising peformance of the ensemble
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    model should form the basis of the last recommendation: optimal training set bias and its interaction with
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    connectance and binary classifier is, in a sense, an hyperparameter that should be assessed. The
    distribution of results in fig. 4 and fig. 5 show that there are variations around the trend; furthermore,
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370
    networks with different structures than the one we simulated here may respond in different ways.
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Figure 1: Consequences of changing the classifier skills (s) and bias (s) for a connectance  $\rho=0.15$ , on accuracy,  $F_1$ , postive predictive value, and  $\kappa$ . Accuracy increases with skill, but also increases when the bias tends towards estimating *fewer* interactions. The  $F_1$  score increases with skill but also increases when the bias tends towards estimating *more* interactions; PPV behaves in the same way. Interestingly,  $\kappa$  responds as expected to skill (being negative whenever s<0.5), and peaks for values of  $b\approx0.5$ ; nevertheless, the value of bias for which  $\kappa$  is maximized in *not* b=0.5, but instead increases with classifier skill. In other words, at equal skill, maximizing  $\kappa$  would lead to select a *more* biased classifier.

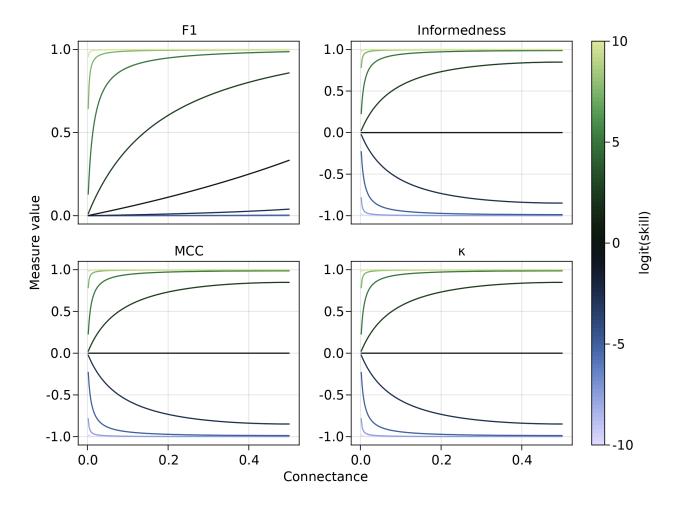


Figure 2: As in fig. 1, consequences of changing connectance for different levels of classifier skill, assuming no classifier bias. Informedness,  $\kappa$ , and MCC do increase with connectance, but only when the classifier is not no-skill; by way of contrast, a more connected network will give a higher  $F_1$  value even with a no-skill classifier.

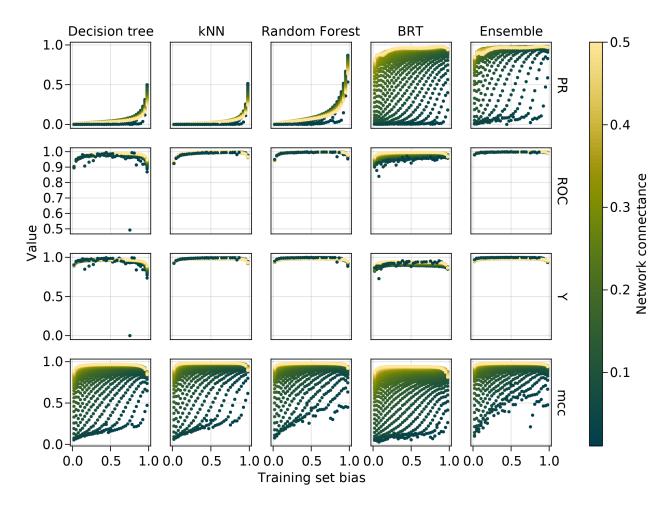


Figure 3: Response of MCC, Informedness, ROC-AUC, and PR-AUC to changes in the training set balance (on the *x* axis) for a series of increasing connectances (color). All of these values approach 1 for a good model, but should be lower when the prediction is more difficult. Informedness is consistently high, and by contrast, MCC increases with additional training set balance. Across all models, training on a more connected network is easier. ROC-AUC is consistently high, and therefore not properly able to separate good from poor classifiers. On the other hand, PR-AUC responds to changes in the training set.

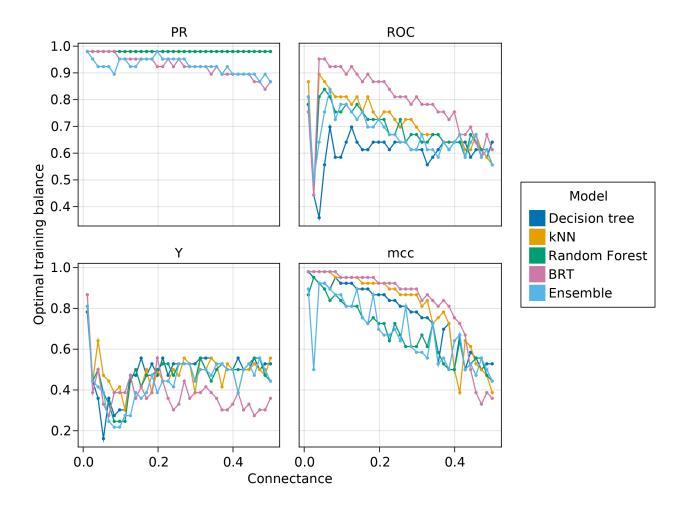


Figure 4: Value of the optimal training set balance for the different models and measures evaluated here, over a range of connectances. Informedness was reliably maximized for balanced training sets, and kept this behavior across models. For other measures, larger connectances in the true network allowed lower biases in the training set. In a large number of cases, "over-correcting" by having training sets with more than half instances representing interactions would maximize the values of the model performance measures.

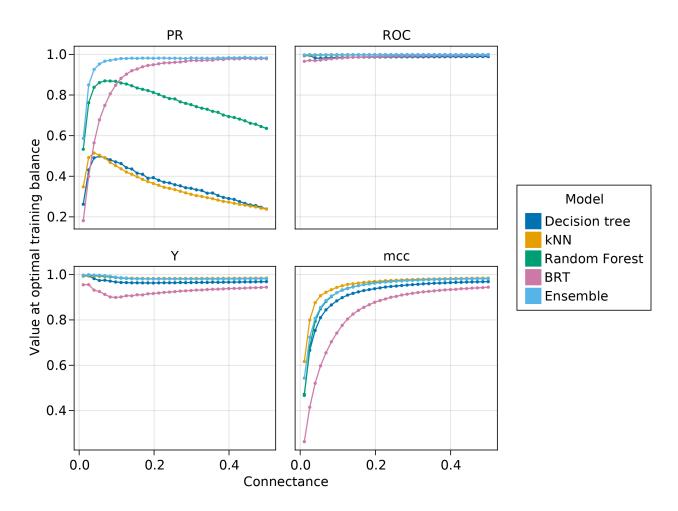


Figure 5: When trained on their optimally biased training set, most models were able to maximize their performance; this is not true for decision tree, which had a very low PR-AUC, and to some extent for ridge regression who had a slow increase with network connectance. The ensemble had a consistently high performance despite incorporating poor models.

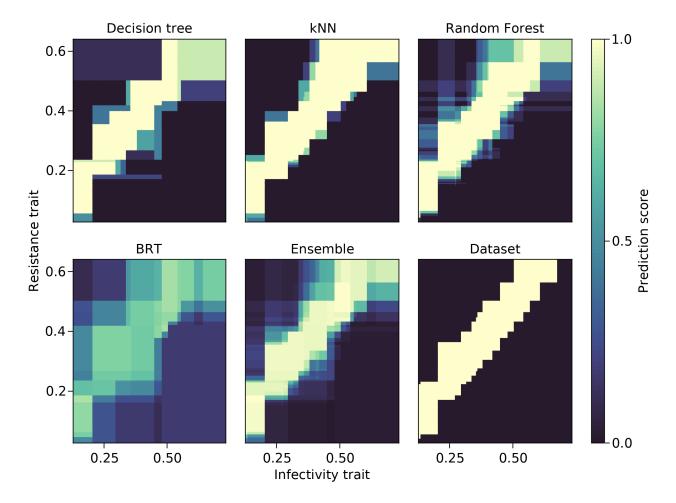


Figure 6: Visualisation of the models predictions for one instance of a network prediction problem (shown in the "Dataset" panel). This figure reveals how inspecting the details of the prediction is important: indeed, although the performance measures hint at the fact that ridge regression is mediocre, this figure reveals that it is making predictions that correspond to a network with an entirely different topology (namely, nested as opposed to diagonal).