

# Guidelines for the supervised learning of species interactions

Timothée Poisot<sup>1,2</sup>

<sup>1</sup> Université de Montréal   <sup>2</sup> Québec Centre for Biodiversity Sciences

## Correspondance to:

Timothée Poisot — [timothee.poisot@umontreal.ca](mailto:timothee.poisot@umontreal.ca)

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1. The prediction of species interaction networks 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.
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. Classifier accuracy and the ROC-AUC are not 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 as a function of the classifier and the network connectance.
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.

- example on diagnostic test: rare events are hard to detect even with really good models
- summary of model challenges for networks
- Strydom et al. (2021) importance of drawing on traits + validation is challenging + comparing across space

Binary classifiers are usually assessed by measuring properties of their confusion matrix, *i.e.* the contingency table reporting true/false positive/negative hits. A confusion matrix is laid out as

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

wherein tp is the number of interactions predicted as positive, tn is the number of non-interactions predicted as negative, fp is the number of non-interactions predicted as positive, and fn is the number of interactions predicted as negative. Almost all measures based on the confusion matrix express rates of error or success as proportions, and therefore the values of these components matter in a *relative* way. 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 focusing on different ways in which the classifier is wrong.

The same approach is used to evaluate *e.g.* species distribution models (SDMs). Indeed, the training and evaluation of SDMs as binary classifiers suffers from the same issue of low prevalence. In a previous work, Allouche et al. (2006) suggested that  $\kappa$  was a better test of model performance than the True Skill Statistic (TSS), which we will refer to as Youden's informedness (or  $J$ ); these conclusions were later criticized by Somodi et al. (2017), who emphasized that informedness' relationship to prevalence depends on assumptions about bias in the model, and therefore recommend the use of  $\kappa$  as a validation of classification performance. Although this work offers recommendations about the comparison of models, it doesn't establish baselines or good practices for training on imbalanced ecological data. Within the context of networks, there are three specific issues that need to be addressed. First, what values of performance measures are we expecting for a classifier that has poor performance? This is particularly important as it can evaluate whether low prevalence can lull us into a false sense of predictive accuracy. Second, independently of the question of model evaluation, is low prevalence an issue for *training*, 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 with the

27 least amount of positive interaction data?

28 In addition to the literature on SDMs, most of the research on machine learning application to life  
29 sciences is focused on genomics (which has very specific challenges, see a recent discussion by Whalen et  
30 al., 2021); this sub-field has generated largely different recommendations. Chicco & Jurman (2020)  
31 suggest using Matthews correlation coefficient (MCC) over  $F_1$ , as a protection against over-inflation of  
32 predicted results; Delgado & Tibau (2019) advocate against the use of Cohen's  $\kappa$ , again in favor of MCC, as  
33 the relative nature of  $\kappa$  means that a worse classifier can be picked over a better one; similarly, Boughorbel  
34 et al. (2017) recommend MCC over other measures of performance for imbalanced data, as it has more  
35 desirable statistical properties. More recently, Chicco et al. (2021) temper the apparent supremacy of the  
36 MCC, by suggesting it should be replaced by Youden's informedness (also known as  $J$ , bookmaker's  
37 accuracy, and the True-Skill Statistic) when the imbalance in the dataset may not be representative  
38 (Jordano, 2016a, which is the case as networks are under-sampled; 2016b), when classifiers need to be  
39 compared across different datasets (for example when predicting a system in space, where undersampling  
40 varies locally; McLeod et al., 2021), and when comparing the results to a no-skill (baseline) classifier is  
41 important. As these conditions are likely to be met with network data, there is a need to evaluate which  
42 measures of classification accuracy respond in a desirable way.

43 A lot of binary classifiers are built by using a regressor (whose task is to guess the value of the interaction,  
44 and can therefore return somethings considered to be a pseudo-probability); in this case, the optimal value  
45 below which predictions are assumed to be negative (*i.e.* the interaction does not exist) can be determined  
46 by picking a threshold maximizing some value on the ROC curve or the PR curve. The area under these  
47 curves (ROC-AUC and PR-AUC henceforth) give ideas on the overall goodness of the classifier. Saito &  
48 Rehmsmeier (2015) established that the ROC-AUC is biased towards over-estimating performance for  
49 imbalanced data; on the contrary, the PR-AUC is able to identify classifiers that are less able to detect  
50 positive interactions correctly, with the additional advantage of having a baseline value equal to  
51 prevalence. Therefore, it is important to assess whether these two measures return different results when  
52 applied to ecological network prediction.

53 We establish that due to the low prevalence of interactions, even poor classifiers applied to food web data  
54 will reach a high accuracy; this is because the measure is dominated by the accidental correct predictions  
55 of negatives. The  $F_1$  score and positive predictive values are less sensitive to bias, but **TODO**

## 56 **Baseline values**

57 Intro

## 58 **Definition of the performance measures**

59  $\kappa$

60  $F_\beta$

61 informedness

62 MCC

63 ROC-AUC

64 PR-AUC

## 65 **Confusion matrix with skill and bias**

66 In this section, we will assume a network of connectance  $\rho$ , *i.e.* having  $\rho S^2$  interactions (where  $S$  is the  
67 species richness), and  $(1 - \rho)S^2$  non-interactions. Therefore, the vector describing the *true* state of the  
68 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  
69 confusion matrix, which ends up expressing *relative* values).

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

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

76 In order to regulate the skill of this classifier, we can define a skill matrix  $\mathbf{S}$  with diagonal elements equal  
 77 to  $s$ , and off-diagonal elements equal to  $(1 - s)$ , and re-express the skill-adjusted confusion matrix as  
 78  $\mathbf{M} \odot \mathbf{S}$ , *i.e.*

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

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

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

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

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

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

85 In all further simulations, the confusion matrix  $\mathbf{C}$  is transformed so that it sums to 1.

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

87 In this section, we will change the values of  $b$ ,  $s$ , and  $\rho$ , and report how the main measures discussed in  
 88 the introduction (MCC,  $F_1$ ,  $\kappa$ , and informedness) are responding to issues with the classifier. Before we do  
 89 so, it is important to explain why we will not focus on accuracy too much. Accuracy is the number of  
 90 correct predictions ( $\text{Tr}(\mathbf{C})$ ) divided by the sum of the confusion matrix. For a no-skill, no-bias classifier,  
 91 accuracy is equal to  $\rho^2 + (1 - \rho)^2$ ; for  $\rho = 0.05$ , this is  $\approx 0.90$ , and for  $\rho = 0.01$ , this is equal to  $\approx 0.98$ . In  
 92 other words, the values of accuracy are expected to be so high that they are not really informative (this is  
 93 simply explained by the fact that 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 of accuracy becomes  $b\rho^2 + (1 - b)(1 - \rho)^2$ , which increases when  $b$  is low, which specifically means that at equal skill, a classifier that under-predicts interactions will have higher accuracy than an un-biased classifier. These issues are absent from balanced accuracy, but should nevertheless lead us to not report accuracy as the primary measure of network prediction success; moving forward, we will focus on other measures.

In order to examine how MCC,  $F_1$ ,  $\kappa$ , and informedness change w.r.t. the imbalance, skill, and bias, we performed a grid exploration of the values of  $\text{logit}(s)$  and  $\text{logit}(b)$  linearly from  $-10$  to  $10$ , of  $\rho$  linearly in  $[0, 0.5]$ , which is within the range of usually observed connectance values for empirical food webs. Note that at this point, there is no food web model to speak of; rather, the confusion matrix we discuss can be obtained for any classification task. Based on the previous discussion, the desirable properties for a measure of classifier success should be: an increase with classifier skill, especially at low bias; a hump-shaped response to bias, especially at high skill, and ideally center around  $\text{logit}(b) = 0$ ; an increase with prevalence up until equiprevalence is reached.

[Figure 1 about here.]

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 bias 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  $\text{logit}(b) = 0$ : 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 towards their correct maximal value when skill changes (*i.e.* a more connected networks will have higher values for a skilled classifier, 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

although sensitive to bias, it is sensitive in a consistent way. Informedness is appropriate at discriminating between different skills, but confounded by bias.  $F_1$  is increasing with bias, and should not be prioritized to evaluate the performance of the model. The discussion of sensitivity to bias should come with a domain-specific caveat: although it is likely that interactions documented in ecological networks are correct, a lot of non-interactions are simply unobserved; as predictive models are used for data-inflation (*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. Furthermore, the weight of positive interactions could be adjusted if some information about the extent of undersampling exists (*e.g.* Branco et al., 2015).

## Numerical experiments on training strategy

In the following section, we will generate random bipartite networks (this works without loss of generality on unipartite networks), and train four binary classifiers (as well as an ensemble model using the sum of ranged outputs from the component models) on 30% of the interaction data. Networks are generated by picking a random infectiousness trait  $v_i$  for 100 species (from a  $B(6, 8)$  distribution), and a resistance trait  $h_j$  for 100 species (from a  $B(2, 8)$  distribution). There is an interaction between  $i$  and  $j$  when  $v_i - \xi/2 \leq h_j \leq v_i + \xi/2$ , 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 fully interval networks that are close analogues to the bacteria–phage model of Weitz et al. (2005), with both a modular structure and a non-uniform degree distribution. This model is easy to learn: when trained with features  $[v_i, h_j, \text{abs}(v_i, h_j)]^T$  to predict the interactions between  $i$  and  $j$ , all four models presented below were able to reach almost perfect predictions all the time (data not presented here) – this is in part because the rule is fixed for all interactions. In order to make the problem more difficult to solve, we use  $[v_i, h_j]$  as a feature vector, and therefore the models will have to uncover that the rule for interaction is  $\text{abs}(v_i, h_j) \leq \xi$ .

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 bias) to be positive, so that the training set has  $\nu n$  interactions, and  $(1 - \nu)n$  non-interactions. We vary  $\nu$  uniformly in  $]0, 1[$ . This allows to evaluate how the measures of binary classification performance respond to artificially rebalanced dataset for a given



150 network connectance. The rest of the dataset ( $n = 7000$  pairs of species) is used as a testing set, on which  
151 all further measures are calculated. Note that although the training set is balanced, the testing set is not,  
152 and retains (part of) the imbalance of the original data.

153 The dataset used for numerical experiments is composed of 64000 such  $(\xi, \nu)$  pairs, on which four  
154 machines are trained: a decision tree regressor, a boosted regression tree, a ridge regressor, and a random  
155 forest regressor. All models were taken from the MLJ.jl package (Blaom et al., 2020; Blaom & Vollmer,  
156 2020) in Julia 1.7 (Bezanson et al., 2017). All machines use the default parameterization; this is an obvious  
157 deviation from best practices, as the hyperparameters of any machine require training before its  
158 application on a real dataset. As we use 64000 such datasets, this would require 256000 unique instances  
159 of tweaking the hyperparameters, which is not realistic. Therefore, we assume that the default  
160 parameterizations are comparable across networks. All machines return a quantitative prediction, usually  
161 (but not necessarily) in  $[0, 1]$ , which is proportional (but not necessarily linearly) to the probability of an  
162 interaction between  $i$  and  $j$ .

163 In order to pick the best adjacency matrix for a given trained machine, we performed a thresholding  
164 approach using 500 steps on predictions from the testing set, and picking the threshold that maximized  
165 Youden's informedness, which is usually the optimized target for imbalanced classification. During the  
166 thresholding step, we measured the area under the receiving-operator characteristic (ROC-AUC) and  
167 precision-recall (PR-AUC) curves, as measures of overall performance over the range of returned values.  
168 We report the ROC-AUC and PR-AUC, as well as a suite of other measures as introduced in the next  
169 section, for the best threshold. The ensemble model was generated by summing the predictions of all  
170 component models on the testing set (ranged in  $[0, 1]$ ), then put through the same thresholding process.  
171 The complete code to run the simulations is given as an appendix; running the final simulation required  
172 4.8 core days (approx. 117 hours).

173 After the simulations were completed, we removed all runs (*i.e.* pairs of  $\xi$  and  $\nu$ ) for which at least one of  
174 the following conditions was met: the accuracy was 0, the true positive or true negative rates were 0, the  
175 connectance was larger than 0.25. This removes both the obviously failed model runs, and the networks  
176 that are more densely connected compared to the connectance of empirical food webs (and are therefore  
177 less difficult to predict, being less imbalanced; preliminary analyses of data with a connectance larger than  
178 3 revealed that all machines reached consistently high performance).

## Effect of training set bias on performance

In fig. 3, we present the response of MCC and informedness to (i) five levels of network connectance and (ii) a gradient of training set bias, for the four component models as well as the ensemble. All models reached a higher performance on more connected networks, and using more biased training sets (with the exception of ridge regression, whose informedness decreased in performance with training set bias). 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 bias, although this increase became less steep with increasing connectance. Interestingly, the ensemble almost always outclassed its component models. In a few cases, both MCC and informedness started decreasing when the training set bias got too close to one, which suggests that it is possible to over-correct the imbalance.

[Figure 3 about here.]

In fig. 4, we present the same information as fig. 3, this time using ROC-AUC and PR-AUC. ROC-AUC is always high, and does not vary with training set bias. On the other hand, PR-AUC shows very strong responses, increasing with training set bias. 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. As in fig. 3, the ensemble outperforms its component models.

[Figure 4 about here.]

Based on the results presented in fig. 3 and fig. 4, 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 bias, and may therefore be more useful at model comparison.

## Required amount of positives to get the best performance

The previous results revealed that the measure of classification performance responds both to the bias in the training set *and* to the connectance of the network; from a practical point of view, assembling a training set requires to withhold positive information, which in ecological networks are very scarce (and

typically more valuable than negatives, on which there is a doubt). For this reason, across all values of connectance, we measured the training set bias that maximized a series of performance measures. When this value is high, the training set needs to skew more positive in order to get a performant model; when this value is about 0.5, the training set needs to be artificially balanced to optimize the model performance. These results are presented in fig. 5.

[Figure 5 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 bias 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.

[Figure 6 about here.]

When trained at their optimal training set bias, performance still had a significant impact on the performance of some machines fig. 6. 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 exemple 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  $S = 50$  species, a connectance of  $\approx 0.16$  ( $\xi = 0.19$ ), and a training set bias of 0.7. The prediction made on the complete dataset is presented in fig. 7. 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.

230

[Figure 7 about here.]

231 The trained models were then thresholded (again by optimising informedness), and their predictions  
 232 transformed back into networks for analysis; specifically, we measured the connectance, nestedness  
 233 (REF), and modularity (REF). The results are presented in tbl. 1. The random forest model is an  
 234 interesting instance here: it produces the network that looks the most like the original dataset, despite  
 235 having a very low PR-AUC, suggesting it hits high recall at the cost of low precision. Although the  
 236 ensemble was about to reach a very high PR-AUC (and a very high ROC-AUC), this did not necessarily  
 237 translate into more accurate reconstructions of the structure of the network. This result bears elaborating.  
 238 Measures of model performance capture how much of the interactions and non-interactions are correctly  
 239 identified. As long as these predictions are not perfect, some interactions will be predicted at the “wrong”  
 240 position in the network; these measures cannot describe the structural effect of these mistakes. On the  
 241 other hand, measures of network structure can have the same value with interactions that fall at  
 242 drastically different positions; this is in part because a lot of these measures covary with connectance, and  
 243 in part because as long as these values are not 0 or their respective maximum, there is a large number of  
 244 network configurations that can have the same value.

Table 1: Values of four performance metrics, and three network structure metrics, for the predictions presented in fig. 7. The values in **bold** indicate the best value for each column (including ties).

Model	MCC	Inf.	ROC-AUC	PR-AUC	Conn.	$\eta$	$Q$
Decision tree	0.83	0.68	0.95	0.15	0.18	0.53	<b>0.49</b>
BRT	0.76	0.89	0.95	0.65	0.22	0.63	0.43
Random Forest	<b>0.89</b>	<b>0.94</b>	<b>0.99</b>	0.41	<b>0.17</b>	<b>0.48</b>	<b>0.49</b>
Ridge Regression	0.67	0.85	0.89	0.38	0.27	1.0	0.26
Ensemble	0.84	0.91	<b>0.99</b>	<b>0.94</b>	0.19	0.54	0.48
Data					0.16	0.45	0.49

## 245 Guidelines for the assesment of network predictive models

246 The results presented here highlight an interesting paradox: larger networks (with lower connectance)  
 247 require more training set bias in order to maximize model performance fig. 5, but are also more difficult to

248 predict according to MCC and PR-AUC fig. 6. This suggests that the task of network prediction will be  
249 difficult regardless of network size: by being limited by the *frequency* of interactions when the network is  
250 large, and by being limited by the *number* of interactions when the network is small. Nevertheless, based  
251 on the simulations and numerical experiments, it is possible to formulate a series of recommendations for  
252 the evaluation of network prediction models.

253 First, because we should have more trust in reported interactions than in reported absences of interactions,  
254 we can draw on previous literature to recommend informedness as a measure to decide on a threshold  
255 (Chicco et al., 2021); this being said, because informedness is insensitive to bias, the model performance is  
256 better evaluated through the use of MCC fig. 3. Because  $F_1$  is monotonously sensitive to classifier bias  
257 fig. 1 and network connectance fig. 2, MCC should be preferred as a measure of model evaluation.

258 Second, because the PR-AUC responds more to network connectance fig. 6 and training set imbalance  
259 fig. 4, it should be used as a measure of model performance over the ROC-AUC. This is not to say that  
260 ROC-AUC should be discarded (in fact, a low ROC-AUC is a sign of an issue with the model), but that its  
261 interpretation should be guided by the PR-AUC value. This again echoes recommendations from other  
262 fields (Saito & Rehmsmeier, 2015).

263 Thirdly, regardless of network connectance *or* measure to evaluate the model performance, as long as the  
264 network connectance is larger than  $\approx 0.1$ , artificially balancing the training set to have equiprevalence will  
265 give the best possible results. This was true for all models.

266 Finally, it is noteworthy that the ensemble model was systematically better than the component models;  
267 even when poor models were included (Random Forest and Decision Tree), the ensemble was able to  
268 leverage the different biases expressed by the models to make an overall more accurate prediction. We do  
269 not expect that ensembles will *always* be better than single models. In a recent multi-model comparison,  
270 Becker et al. (2021) found that the ensemble was *not* the best model. There is no general conclusion to  
271 draw from this besides reinforcing the need to be pragmatic about which models should be included in the  
272 ensemble, or whether to use an ensemble at all. In a sense, the surprising performance of the ensemble  
273 model should form the basis of the last recommendation: optimal training set bias and its interaction with  
274 connectance and binary classifier is, in a sense, an hyperparameter that should be assessed. The  
275 distribution of results in fig. 5 and fig. 6 show that there are variations around the trend; furthermore,  
276 networks with different structures than the one we simulated here may respond in different ways.

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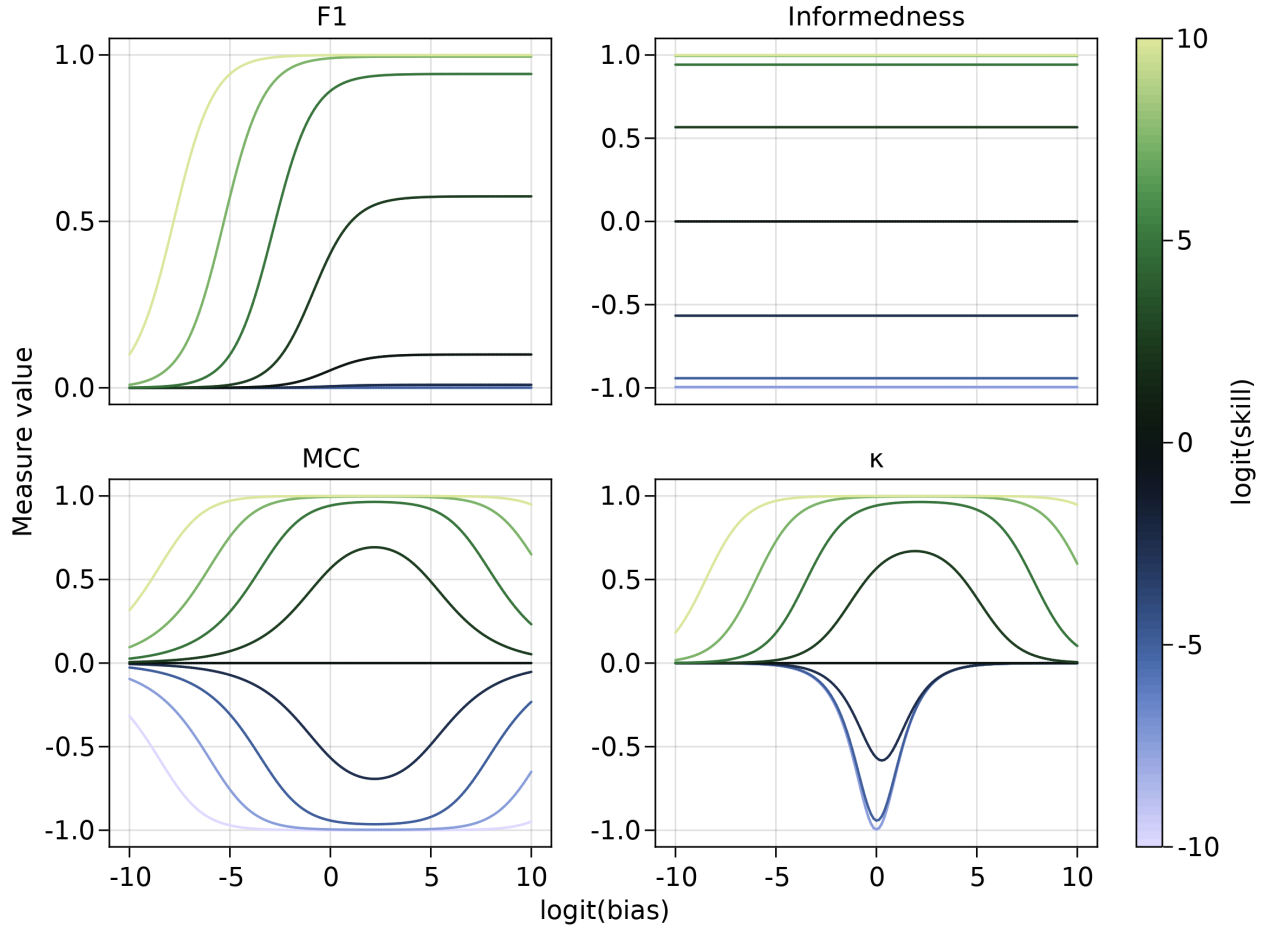


Figure 1: Consequences of changing the classifier skills ( $s$ ) and bias ( $b$ ) for a connectance  $\rho = 0.15$ , on accuracy,  $F_1$ , positive predictive value, and  $\kappa$ . Accuracy increases with skill, but also increases when the bias tends towards estimating *fewer* interactions. The  $F_1$  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  $b \approx 0.5$ ; nevertheless, the value of bias for which  $\kappa$  is maximized is *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.

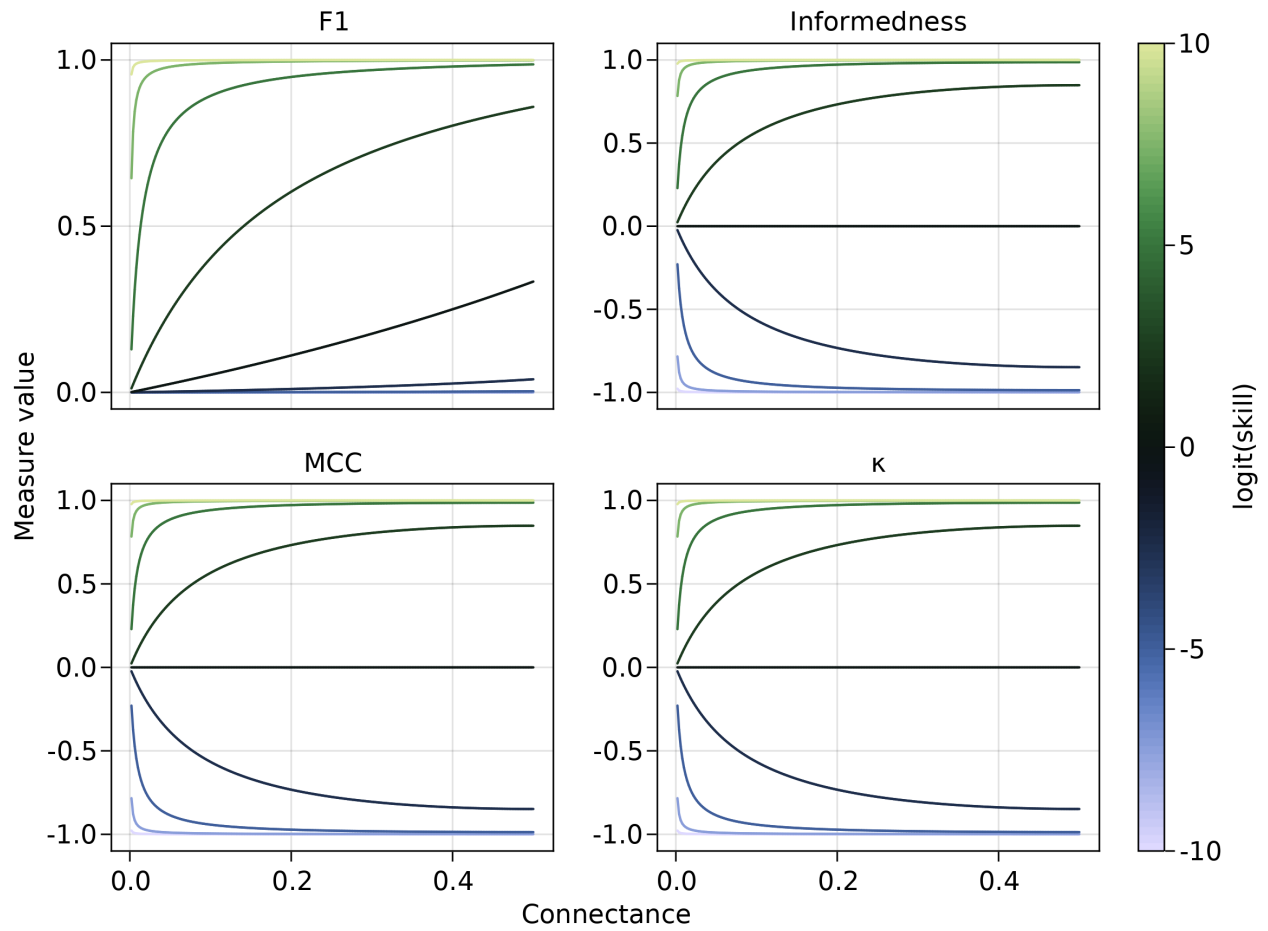


Figure 2: As in fig. 1, consequences of changing connectance for different levels of classifier skill, assuming no classifier bias. Informedness,  $\kappa$ , and MCC do increase with connectance, but only when the classifier is not no-skill; by way of contrast, a more connected network will give a higher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

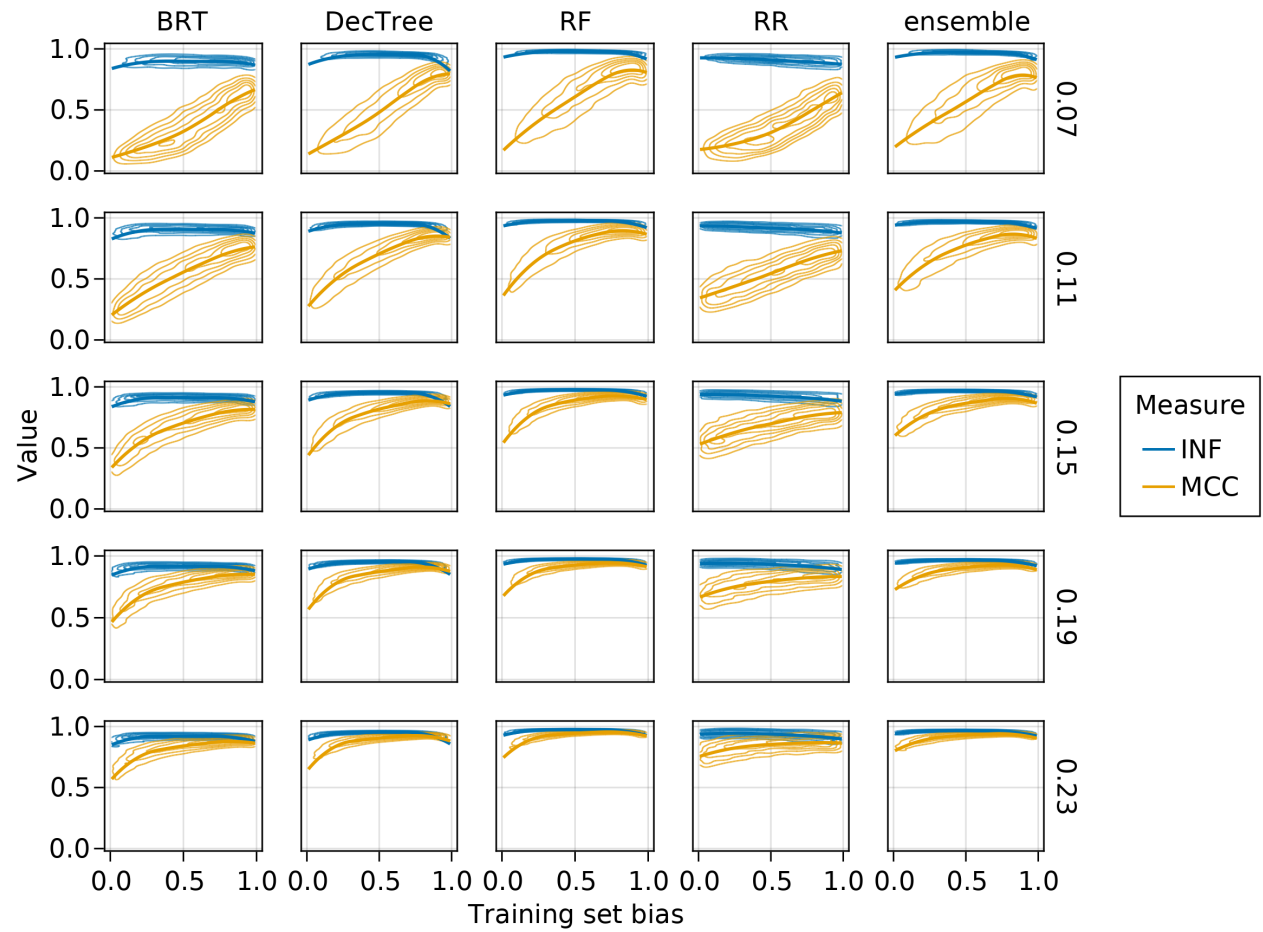


Figure 3: TODO

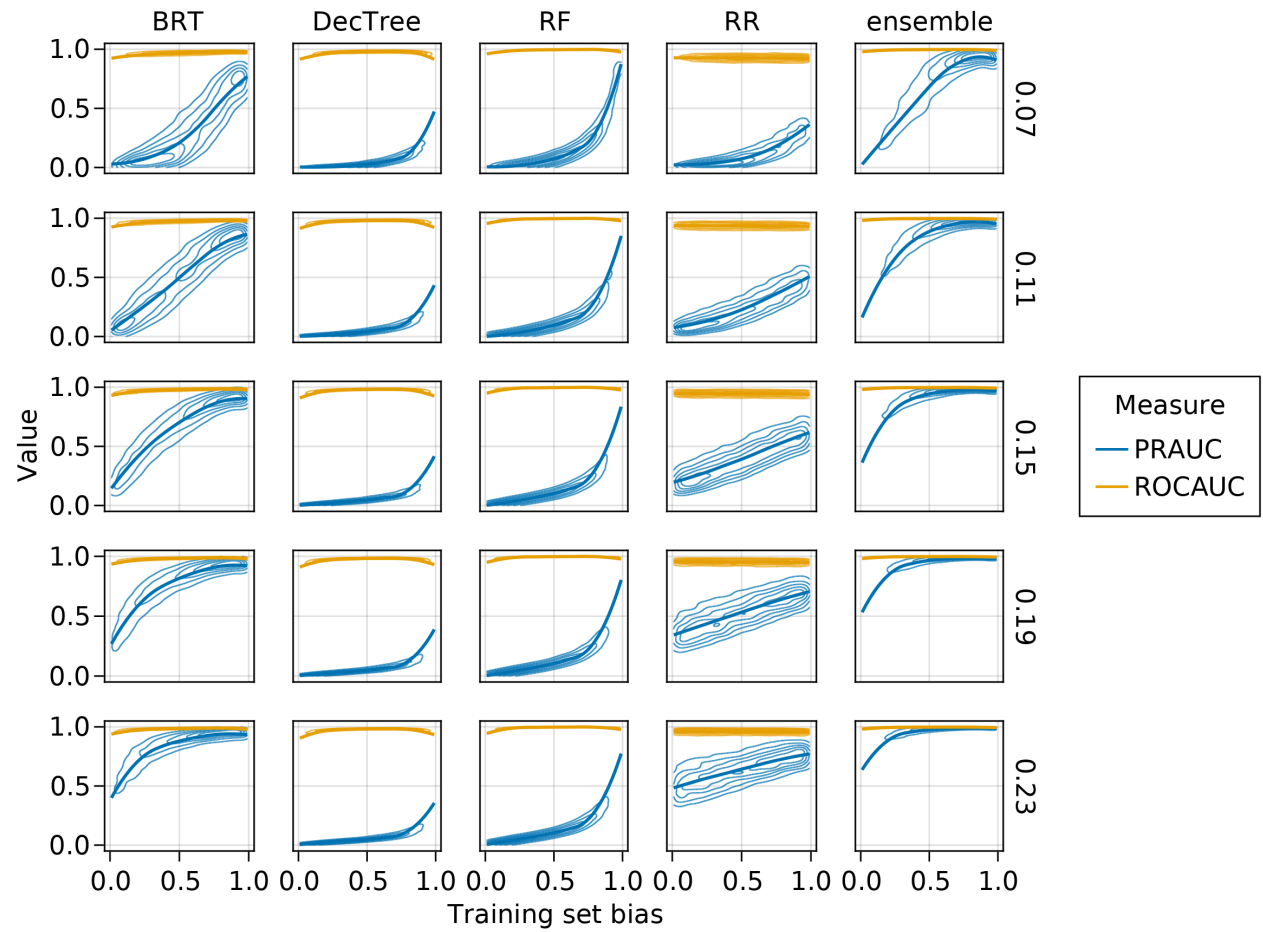


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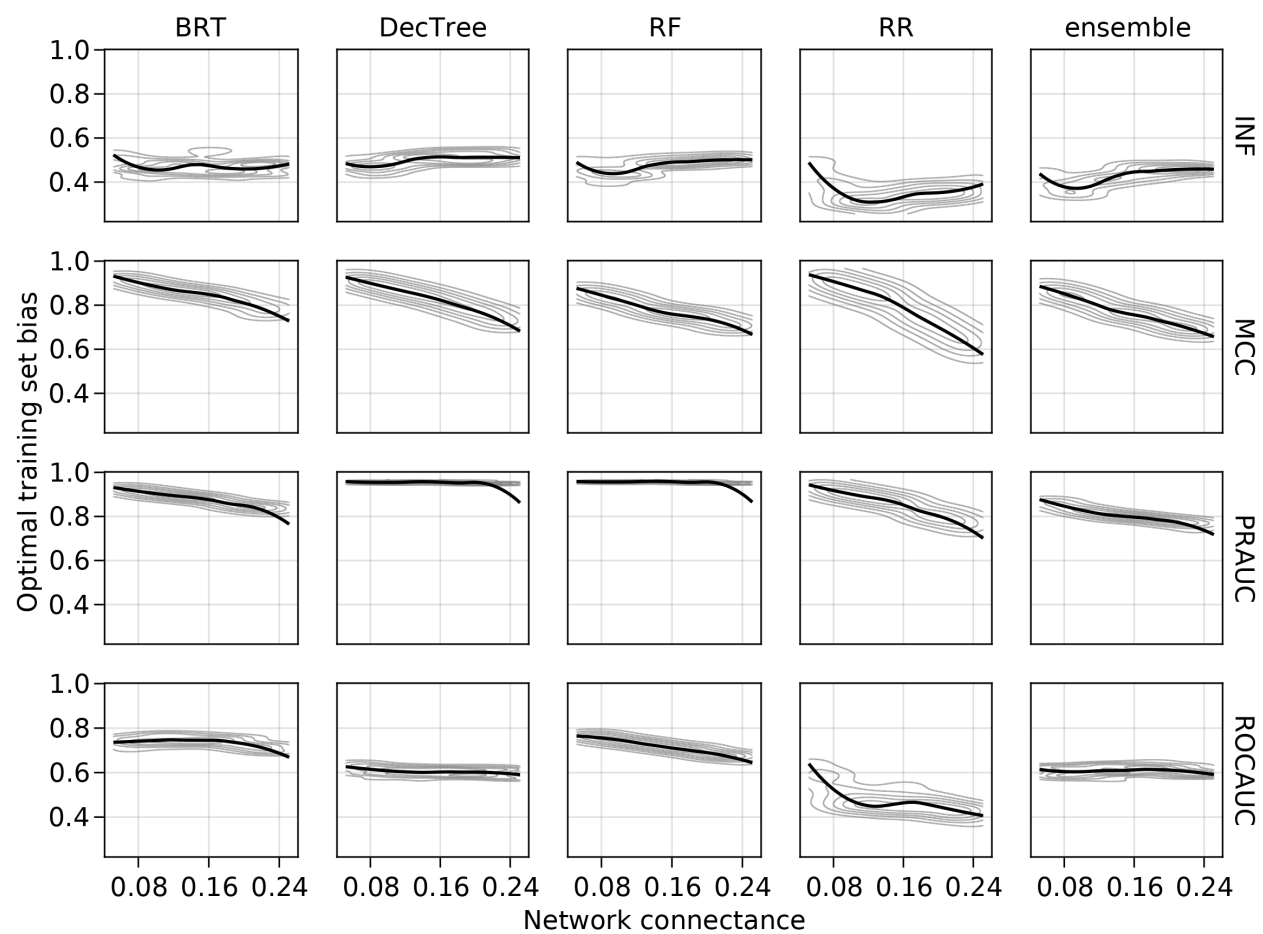


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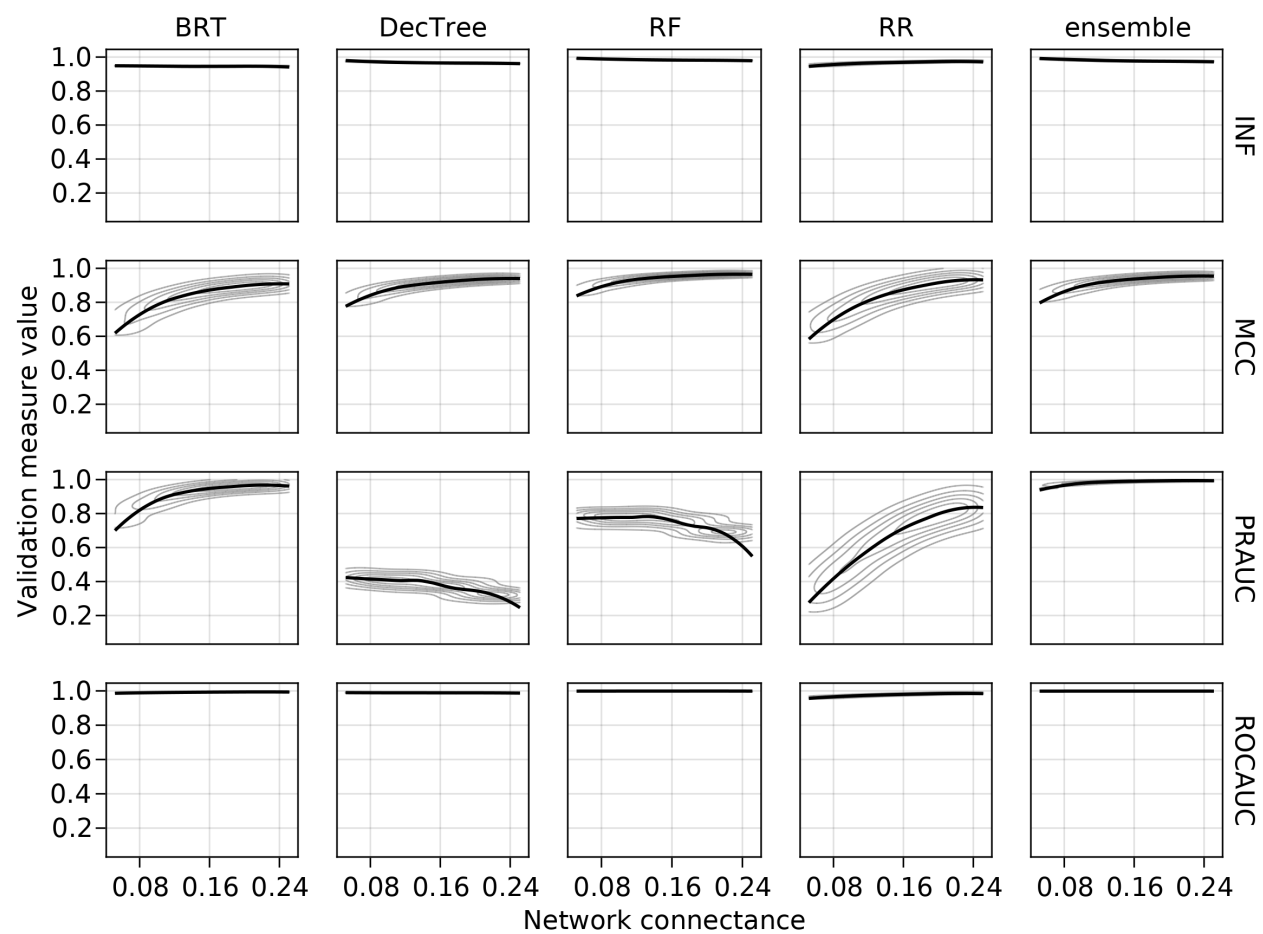


Figure 6: TODO

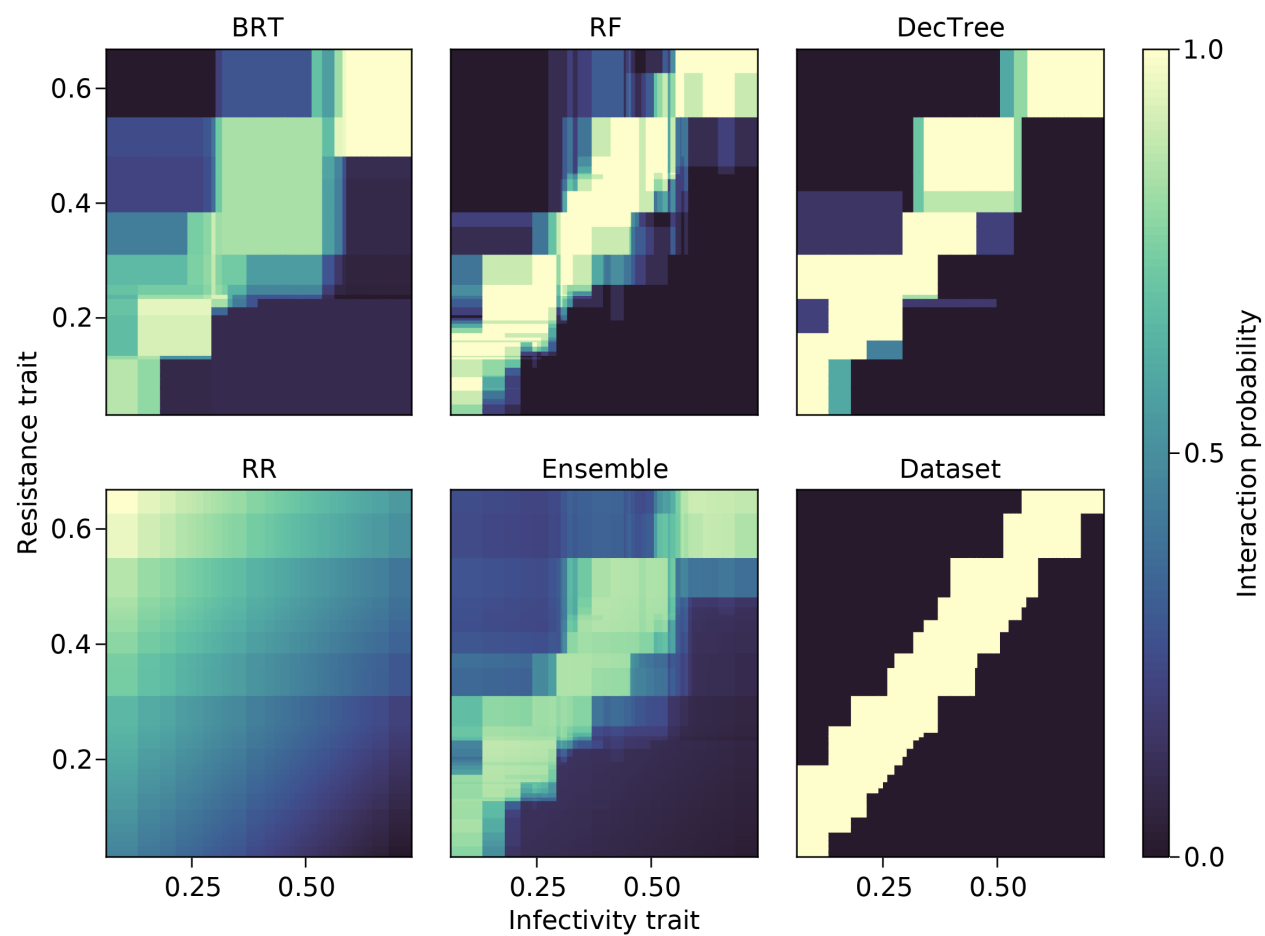


Figure 7: TODO