
Experimental Procedures for Exploiting Structure in AutoML Loss Landscapes

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Abstract Recent observations regarding the structural simplicity of algorithm configuration landscapes have spurred the development of new configurators that obtain provably and empirically better performance. Inspired by these observations, we recently performed a similar analysis of AutoML loss landscapes – that is, the relationship between hyper-parameter configurations and machine learning model performance. In this study, we propose two new variations of an existing, state-of-the-art hyper-parameter configuration procedure. We designed each method to exploit a specific property that we observed common among most AutoML loss landscapes; however, we demonstrate that neither are competitive with existing baselines. In light of this result, we construct artificial AutoML scenarios that allow us to show when the two new methods can be expected to outperform their baselines and when they cannot, thereby providing additional insights into AutoML loss landscapes.

1 Introduction

In recent years, a large variety of methods for automating machine learning have emerged to automatically select a machine learning model and training algorithm along with configuring its hyper-parameters (Bergstra et al., 2011; Bergstra and Bengio, 2012; Jamieson and Talwalkar, 2016; Li et al., 2017, 2020; Fusi et al., 2018; Falkner et al., 2018; Yang et al., 2019), thereby making machine learning faster and easier to use for data scientists and non-experts alike. In parallel to this work, a growing amount of attention has been devoted to studying the landscapes – that is, the relationship between algorithm parameter settings and performance – of a related problem: The configuration of algorithms for \mathcal{NP} -hard problems (Pedersen, 2010; Yuan et al., 2010, 2012; Pushak and Hoos, 2018; Harrison et al., 2019, 2020, 2021; Asi and Duchi, 2019; Treimun-Costa et al., 2020; Cleghorn and Ochoa, 2021). These insights – for example, the observation that individual algorithm parameters tend to elicit uni-modal responses in performance (Pushak and Hoos, 2018) – have already successfully translated into new, state-of-the-art configurators with provably (Hall et al., 2020) and empirically (Pushak and Hoos, 2020) better performance.

Comparatively little is known about AutoML loss landscapes (Garciaarena et al., 2018; Pimenta et al., 2020), which relate the hyper-parameter settings of a given ML algorithm to the validation loss achieved by it. This gap inspired our own recent study of AutoML loss landscapes (Pushak and Hoos, view), in which we observed two common trends in the hyper-parameters of individual machine learning models over a range of widely-studied benchmarks: First, they appear to induce landscapes that are uni-modal (or very close to it); and second, they appear to have relatively benign hyper-parameter interactions. In particular, we observed that a simplistic configuration procedure that optimizes each hyper-parameter a single time, in a random order, can find hyper-parameter configurations tied with optimal with very high probability. In light of these observations, existing hyper-parameter configuration procedures may be spending larger-than-necessary amounts of

their budgets evaluating configurations in regions of the hyper-parameter configuration space that can be reasonably assumed to be of poor quality.

For this reason, using relatively simple surrogate models that incorporate our prior expectation regarding the structure of the landscapes may help to improve the efficiency of configuration procedures by cutting down on unnecessary configuration evaluations. However, both of our proposed surrogate models (see Section 2) proved to be too simple, resulting in similar or worse performance compared to BOHB (Falkner et al., 2018) for most of the real-world AutoML scenarios we studied (see Section 4). To investigate why, we evaluated the methods on four hand-crafted benchmark scenarios (see Section 3), and thus obtain additional insights into a) the structure of AutoML loss landscapes, b) the reasons why these particular methods did not perform as well as expected, and c) the conditions under which they can be expected to perform well (see Section 5).

2 Experimental AutoML Procedures

We propose two modifications to BOHB (Falkner et al., 2018) (which uses Bayesian optimization) that both involve replacing the tree-structured Parzen estimator (TPE) (Bergstra et al., 2011) model with simpler, more constrained surrogate models chosen to reflect one of the structural properties that we observed in AutoML loss landscapes (Pushak and Hoos, view):

- **Convex quadratic approximation (CQA) surrogate models** (Rosen and Marcia, 2004). In fact, the results of our landscape analysis revealed that AutoML loss landscapes are not convex; however, it may still be helpful to use a convex quadratic function as a surrogate model for the landscape, as it is perhaps the simplest n -dimensional model that is constrained to be uni-modal.
- **B-spline basis surrogate models.** Given that interactions between hyper-parameters in AutoML loss landscapes are relatively benign, fitting a linear model to a B-spline basis function (Eilers and Marx, 1996) is a natural choice. These models are known to provide stable approximations (Bickel et al., 2009) when features do not interact, because they force the fitted model to contain no (hyper-parameter) interactions.

Similar to BOHB’s TPE model, each of our surrogate models attempt to approximate the full AutoML loss landscapes. In the case of the convex quadratic surrogate model, this resulted in substantially increased overhead for fitting each of the models. Therefore, in order to avoid slowing down the configuration procedure unnecessarily, we fit the models asynchronously. However, this means that we are required to suggest multiple configurations using the same surrogate model. As a result, we employed a similar, but adapted version of BOHB’s method for suggesting configurations, which does not completely optimize the surrogate model. For more details about each of these design decisions as well as how we fit each model, see Appendix A.

3 Experimental Setup

The primary goal in this study is to answer the question: “Can AutoML loss landscape structure be exploited to improve configurator performance?” Therefore, the strategy used to allocate budget between the evaluation of each candidate configuration is effectively an orthogonal design decision that acts as a confounding factor. To isolate the impact of the search method, our preliminary analysis involved only two baselines: BOHB (Falkner et al., 2018) and Hyperband (Li et al., 2017). To further eliminate confounding factors, we used the implementation of Hyperband that is available with the original implementation of BOHB, and used the same values for all of the shared parameters for all four of the configurators.

We evaluated each configurator on five scenarios, spanning a range of different machine learning methods and datasets, as well as on four hand-crafted benchmark scenarios. We present

a summary of the scenarios in Table 2 in Appendix B.¹ For each of the real-world scenarios, we performed 25 independent runs of the configurators, and we report the median test loss with 95% bootstrap percentile confidence intervals. To obtain a stable estimate for the test loss, we performed 5 independent training runs of the machine learning methods using each configuration and recorded their mean test losses. For the hand-crafted benchmark scenarios, which require substantially less running time to use, we performed 101 independent runs of the configurators and we report the exact loss underlying the landscape.

We also evaluated the four configurators on four artificial scenarios that simulate a binary classifier’s error rate, p , such that $0 < p < 1$, thereby allowing us to control the precise shape of the AutoML loss landscape optimized by the configuration procedures. This control allowed us to test hypotheses about the relative behaviour of the methods under various circumstances, thereby providing answers to questions that arose from our results on the real-world benchmark scenarios. We motivate and define the precise landscapes that we used for the simulated classifier scenarios in Section 4, based on the results of the real-world scenarios.

4 Results

We show the anytime configuration results for three configuration budgets for the five real-world scenarios in Table 3 in Appendix C. Overall, none of the methods consistently perform substantially better than any of the others; however, BOHB found the best or tied for the best on each scenario for all configuration budgets. Unfortunately, while both of the new methods we proposed can find high-quality configurations, neither of them appear to consistently work as well as BOHB.

Why does a CQA surrogate model not work better? Given that we observed that most AutoML loss landscapes are uni-modal (Pushak and Hoos, view), it seems reasonable to assume that a uni-modal surrogate function should be able to improve the performance of BOHB. However, since the CQA surrogate model is a quadratic function, it implicitly assumes that the landscape is symmetric around the globally optimal configuration – which was frequently not the case in our previous study. Therefore, this could be at least partly responsible for the poor performance that we observed with the CQA surrogate model.

To investigate this hypothesis, we compared the performance of the four configuration procedures on two artificial scenarios, where the machine learning algorithm’s loss is simulated as described in Appendix B. In both cases, the simulated classifier had a single real-valued hyperparameter, x , that was searched over the range of values $[-1, 1]$. For the first scenario, we set the error rate to be

$$p_{\text{symmetric}}(x) = |x|^3 + 0.01 \tag{1}$$

and for the second scenario we set it to

$$p_{\text{asymmetric}}(x) = \begin{cases} |x|^3 + 0.01 & \text{if } x < 0 \\ \frac{1}{5} \cdot |x|^3 + 0.01 & \text{otherwise.} \end{cases} \tag{2}$$

In the asymmetric scenario, it *should* be easier for the configuration procedures to find better-quality configurations, because 50% of the configurations obtain better error rates. However, we observed that when using a CQA surrogate model, the anytime configuration performance on the asymmetric scenario was actually worse than on the symmetric scenario (see Table 1). Furthermore, we see that on the symmetric scenario, using the CQA surrogate model yields significantly better

¹We were unable to reproduce qualitatively similar results for BOHB (Falkner et al., 2018) and Hyperband (Li et al., 2017) on two of the benchmark scenarios we studied (see Appendix B). However, from personal communication with the authors, we confirmed that we correctly set up the experiments and reproduced their method of analysis. To the best of our knowledge, the only differences in our execution environments were the particular machines used and, perhaps, the versions of some of the software packages.

Table 1: Results from applying the hyper-parameter configurators to the artificial, hand-crafted AutoML scenarios. At each configuration budget (the cumulative number of training examples used to “train” candidate models) we show the median loss (percentage of errors) over the independent configurator runs with 95% bootstrap percentile confidence intervals. The median losses are shown in boldface if they are within 0.01% of the best loss for a given configuration budget. Smaller is better.

	Symmetric	Asymmetric	No Interactions	Interactions
Budget (10%)	13k Examples	13k Examples	13k Examples	13k Examples
CQA	1.01 [1.01, 1.02]	1.04 [1.03, 1.09]	3.56 [2.89, 4.88]	3.08 [2.73, 3.70]
Spline	1.06 [1.03, 1.10]	1.08 [1.04, 1.13]	5.19 [3.93, 6.24]	3.50 [2.99, 4.19]
Random (HB)	1.11 [1.05, 1.20]	1.08 [1.05, 1.12]	5.26 [3.75, 6.39]	4.12 [3.40, 4.43]
TPE (BOHB)	1.12 [1.07, 1.25]	1.08 [1.04, 1.14]	4.32 [3.36, 5.48]	3.68 [3.03, 4.32]
Budget (50%)	67k Examples	67k Examples	67k Examples	67k Examples
CQA	1.01 [1.00, 1.01]	1.02 [1.01, 1.03]	2.12 [1.86, 2.29]	1.27 [1.19, 1.37]
Spline	1.02 [1.01, 1.04]	1.02 [1.01, 1.03]	1.27 [1.20, 1.44]	1.60 [1.47, 1.74]
Random (HB)	1.04 [1.02, 1.05]	1.02 [1.01, 1.03]	2.06 [1.84, 2.40]	1.91 [1.77, 2.25]
TPE (BOHB)	1.04 [1.03, 1.10]	1.02 [1.01, 1.04]	2.40 [1.90, 2.83]	1.64 [1.38, 2.04]
Budget (100%)	135k Examples	135k Examples	135k Examples	135k Examples
CQA	1.00 [1.00, 1.00]	1.01 [1.00, 1.01]	1.32 [1.21, 1.47]	1.15 [1.12, 1.18]
Spline	1.01 [1.00, 1.02]	1.01 [1.00, 1.02]	1.11 [1.09, 1.13]	1.26 [1.19, 1.31]
Random (HB)	1.02 [1.01, 1.04]	1.01 [1.01, 1.02]	1.65 [1.54, 1.85]	1.59 [1.48, 1.73]
TPE (BOHB)	1.03 [1.01, 1.04]	1.01 [1.01, 1.02]	1.38 [1.23, 1.59]	1.27 [1.16, 1.32]

performance than all of the other methods for most configuration budgets; however, on the asymmetric scenario it instead ties with all of the other methods for all but the 10% configuration budget, at which point it still has a small advantage.

This example clearly illustrates a rather large failure mode for using a CQA model as a surrogate function; even though the asymmetric artificial classifier scenario aligns very closely with the types of landscapes for which we had designed it to perform well, we see that it does not outperform any of the other methods at all.

Why does the spline surrogate model not work better? Similar to the CQA model, we had expected the spline surrogate model to yield better results, given that we observed that most hyper-parameters could be safely configured independently in a random sequence (Pushak and Hoos, view). However, note that this definition of simplicity among hyper-parameter interactions may not actually indicate that a *surrogate model* can assume that the hyper-parameters are independent of each other. Indeed, the function $f(x, y) = |x - y|$ is trivial to minimize when considering either x or y independently – for any fixed value of one, the other can be adjusted to yield an optimal solution. Despite this “simplicity”, a the spline surrogate model would be unable to accurately model the function due to the strong interaction between the hyper-parameters x and y .

To investigate, we again compared the performance of the four configuration procedures on two artificial classifier scenarios. Each scenario contained two hyper-parameters, x and y in $[-1, 1]$. In one case there were no interactions and the error rate was defined as

$$p_{\text{no interactions}}(x, y) = \frac{1}{2} \cdot |x| + 0.01 \quad (3)$$

(y was ignored). The second case used the same objective function after applying a 45 degree rotation, that is, the error rate was set to

$$p_{\text{interactions}}(x, y) = \frac{1}{2 \cdot \sqrt{2}} \cdot |x - y| + 0.01. \quad (4)$$

In this experiment, the scenario with hyper-parameter interactions *should* be easier for the configurators, because the basin of optimal solutions lies along the diagonal $x = -y$ (which has length $2 \cdot \sqrt{2}$ for $x, y \in [-1, 1]$) instead of along the line $x = 0$ (which has length 2 over the same domain). However, we see that this simple rotation of the landscape causes the scenario to become substantially more challenging for the spline model. For both of the larger two configuration budgets, the spline model finds configurations with smaller error rates when there are no hyper-parameter interactions (see Table 1). Similarly, for all three configuration budgets, we see that the spline model finds the best solutions when there are no interactions, whereas when there are interactions, the CQA model performs best. (Note that both of the landscapes in these scenarios are perfectly symmetric along two axes around the global optima, hence we would expect CQA to do well in both cases – which it does.)

This example clearly illustrates an important failure mode for the spline model; even though it may often be sufficient to configure each hyper-parameter independently in a random sequence, it is not necessarily safe to build surrogate models that assume that they are independent.

Furthermore, similar to the parameters of the meta-heuristics studied by Yuan et al. (2012), we expect there to be some hyper-parameters with this type of compensatory interaction in AutoML scenarios. For example, consider the learning rate (`l_rate`) and the momentum decay rate (`mdecay`) hyper-parameters for the Bayesian neural network scenarios. The learning rate corresponds to the step size taken in each iteration of the optimization process, and the momentum decay rate corresponds to how quickly old gradient information is lost. Specifically, a large value for momentum corresponds to making heavy use of the previous gradients, whereas a small value corresponds to heavily trusting the most recently observed gradient.

In cases where gradient measurements are noisy, it is helpful to use large values of momentum to take an average over a larger number of previous gradient observations, thereby smoothing the path taken by the optimizer. This should be particularly important for large step sizes, since otherwise the optimizer will be taking large steps in nearly-random directions, effectively placing too much trust in each gradient observation. Indeed, we empirically observed precisely this relationship between these two hyper-parameters. For example, the best 5% of the 1 234 anytime configurations suggested by any of the configuration procedures have a correlation coefficient of 0.83 for the `l_rate` and `mdecay` hyper-parameters on the protein structure dataset. These observations likely explain why the spline model did not produce better results on the real-world scenarios.

5 Conclusions and Future Work

We introduced two experimental modifications to a state-of-the-art AutoML hyper-parameter configuration procedure, BOHB (Falkner et al., 2018). Both of the modifications we made were inspired by the simplicity of the structure we observed in most AutoML loss landscapes (Pushak and Hoos, view). However, in both cases, the models were too simple to be effective on real-world problems, which has led to a deeper understanding of the structure of AutoML loss landscapes.

In our first attempt, we replaced the TPE model in BOHB with a convex quadratic approximation (CQA) surrogate model. We assumed that because the CQA model is uni-modal (like most AutoML loss landscapes), it should help guide the search process away from regions of the configuration space that we can reasonably assume contain poor configurations. However, the CQA model implicitly assumes that the landscape is symmetric around the global optimum, which proved to be too strong of an assumption.

Our second attempt used B-spline models, which are known to provide high-quality approximations of functions with no interactions between the variables. This choice was motivated by our observation that configuring each hyper-parameter independently, a single time and in a random order, yields configurations tied with optimal with very high probability (Pushak and Hoos, view). However, this can still occur in the presence of strong *compensatory hyper-parameter interactions* – that is, when the basin of optimal solutions is diagonally oriented within the landscape. In practice,

we expect this type of interaction to occur between some hyper-parameters. For example, we showed that the high-quality configurations for the Bayesian neural network scenarios have a correlation coefficient of 0.83 between the values of their momentum decay rate and learning rate hyper-parameters.

Our results suggest a natural next question: What surrogate model can be used to encode our priors, without being too simple, thereby making assumptions that are too strong?

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A Extended Methods

In this section, we include additional details about our two experimental hyper-parameter configuration methods. Each method was implemented as a modification to the open-source version of BOHB (Falkner et al., 2018), which has been released with a BSD 3-clause license at <https://github.com/automl/HpBandSter>.

A.1 Convex Quadratic Surrogate Models

To fit convex quadratic models, we use the methods proposed by Rosen and Marcia (2004). One advantage of their methods compared to others is that the models are fitted using the L_1 norm, which is more robust than the L_2 norm in the presence of noise and outliers. They further impose the constraint that the fitted models must under-estimate all of the available training data. While it is unclear whether or not this will be beneficial for our application, we do not expect it to harm the performance of the method. Adding such a constraint will increase the residuals of an optimally-fitted model and thus will decrease the accuracy of the model’s predictions. However, a model need not have high predictive accuracy in order to be an effective surrogate model; indeed, an ideal surrogate model only needs two properties: it should be easy to optimize and its global optimum should be in a similar location to the global optimum of the original landscape.

To fit the convex quadratic under-estimator models, Rosen and Marcia (2004) proposed two methods, both of which formulate the problem as a two-step procedure. In the first step of each, the problem is simplified by imposing additional constraints on the quadratic model, thereby allowing the problem to be efficiently solved using linear programming. They then proposed to use the incumbent solution from the linear program to initialize the parameters of the non-linear program that is used to fit the general convex quadratic approximation model. This initial guess for the solution to the non-linear program can then be iteratively refined using an interior-point method.

In the first method, they force the Hessian to be a diagonal matrix, and in the second, they force it to be a diagonally-dominant matrix. Rosen and Marcia (2004) found that the second method yielded higher-quality final solutions, therefore, we chose to use it when possible. For a set of m data points in \mathbb{R}^n , the general convex quadratic model fit using this procedure uses $s = (n + 1) \cdot (n + 2)/2$ variables, and thus requires at least $m \geq s$ data points before models can be fit. However, the separable convex quadratic model obtained using a diagonal Hessian matrix only uses $s = 2 \cdot n + 1$ variables. We therefore fit a separable convex quadratic surrogate model whenever $2 \cdot n + 1 < m < (n + 1) \cdot (n + 2)/2$, and we fit a general convex quadratic surrogate model when $m \geq (n + 1) \cdot (n + 2)/2$.

The methods proposed by Rosen and Marcia (2004) also require a parameter ϵ , which is used to constrain the minimum eigenvalue of the Hessian, thereby improving the conditioning of the surrogate-fitting optimization problems and ensuring that the optimum of the surrogate is unique. In our experiments, we set $\epsilon = 10^{-3}$, since it is well-known that some hyper-parameters may have relatively little impact on the solution quality (see for example, Hutter et al. (2014)). Furthermore, we optimized the linear programs using the default linear-programming solver available in python’s SciPy package (Virtanen et al., 2020). To optimize the non-linear programs, we used SciPy’s sequential least squares quadratic programming solver (SLSQP).

A.2 B-Spline Basis Surrogate Models

Gaussian processes are a more commonly used model in AutoML methods that would likely provide similar-quality results to spline models. Gaussian processes are typically chosen, because they can provide a confidence interval around their predictions, which is necessary for most Bayesian optimization methods. However, in our application, where we wish to focus on exploitation rather than trading off between exploration and exploitation, we do not require this confidence interval. Instead, we prefer splines for their lower computational complexity.

Another advantage of using splines is that they support various methods for extrapolation. For our application, we chose to extrapolate using a constant function, as this will ensure that unpromising directions for exploration remain unexplored, whereas promising directions for exploration will continue to be explored.

To compute the spline basis function, we used the implementation available in `scikit-learn`'s latest development branch,² with the default settings of its parameters: five knots per feature with a third-degree basis. For hyper-parameters that are searched on a linear scale, we used a uniform spacing for the knots, whereas we spaced the knots geometrically for hyper-parameters that are searched on a logarithmic scale. Because we anticipated that the models would frequently be under-constrained, particularly early in the configuration process when relatively little configuration performance data is available, we used LASSO regression with five-fold cross-validation to choose the value of the regularization parameter.

A.3 Asynchronous Model Fitting and Selection

Since fitting these surrogate models typically requires substantially more time than fitting a TPE model, we fit the models asynchronously. In BOHB, new models are fitted when worker processes report new configuration results to the main process. In our methods, we initiate and check for completed model-fitting processes in this function as well (see Algorithm 1). We limit the number of models being fitted at any given time to one, to ensure that it does not overburden the machine and slow down the evaluation of the configurations.

Algorithm 1 The report function for our experimental AutoML configurators.

```

1: input
2:    $m_I(c)$ , the performance of a newly evaluated configuration
3:    $f$ , the level of fidelity at which the configuration was evaluated
4:    $R$ , a dictionary mapping fidelities to configuration evaluation results
5:    $model$ , a dictionary mapping fidelities to previously-fit surrogate models (or None)
6: output
7:   None
8: procedure report
9:   # Save the configuration results and check for model-fitting results
10:   $R[f][c] := m_I(c)$ 
11:  Check if a model-fitting process has completed
12:  # Save the new model, if applicable
13:  if a model-fitting process reported  $model_{new}$  at fidelity level  $f'$ , then
14:    if  $model[f']$  is None or  $model_{new}$  better approximates  $R[f']$  than  $model[f']$ , then
15:       $model[f'] := model_{new}$ 
16:  # Start a new model-fitting process, if applicable
17:  if no model-fitting processes are running, then
18:    for each fidelity  $f \in R$  in descending order, do
19:      if  $length(R[f]) \geq 2 \cdot |P| + 1$ , then
20:        Initiate a model-fitting process with  $R[f]$ 
21:      break
22:  return

```

Similarly, BOHB also evaluates each candidate configuration asynchronously. Each time a worker process for BOHB is done evaluating a configuration, the worker process reports its results to the main process. When the main process receives this data, it checks to see if there is sufficient data available to fit a model and if there are no models that are currently being fitted. If both of these conditions are met, the main process will initiate an asynchronous model-fitting process. Once the model-fitting process is complete, it reports its results back to the main process.

²As of 2021-06-24.

Similar to BOHB, we fit separate CQA or spline models for each level of fidelity of the performance estimates. In some early prototypes, we experimented with using a weighted combination of some or all of the available performance estimates, regardless of their level of fidelity. However, we observed that because some measures of fidelity not only decrease the variability in the performance estimates, but also substantially change the performance values for a given configuration, doing so typically caused the model-fitting procedure to diverge, thereby producing unreliable surrogate models. Therefore, as soon as sufficient data to fit a model is available for a higher-fidelity performance estimate, the methods switch to fitting models using the higher level of fidelity instead.

In some rare cases the addition of new configuration performance data, or the switch from a large training set with low-fidelity performance estimates to a small training set with higher-fidelity performance estimates, can de-stabilize the model-fitting procedure, thereby causing it to diverge. To protect against this, we only accept a new surrogate model if the mean absolute error of a newly-fitted model is not worse than the current accepted model, if one exists.

A.4 Suggesting a Configuration

The simplest way to use a surrogate model to guide an optimization process is to find the globally optimal solution to the surrogate model and then to evaluate that solution using the original objective function. When the surrogate model is a convex quadratic model, this would be both easy and inexpensive to do. However, surrogate models that are fitted to previous performance data are typically only good approximations of the original landscape near to the previously-evaluated solutions. Therefore, it is common practice to employ, for example, a trust region, which limits how far away from the current incumbent the surrogate model should be trusted. The next configuration to evaluate is then the configuration predicted to have the optimal performance by the surrogate model that is contained within the trust region.

However, in our application, the surrogate models are fitted asynchronously, so it is also sometimes necessary for multiple new configurations to be suggested using the same surrogate model. Therefore, rather than using a trust region, we employ a somewhat similar trick to BOHB, whereby we randomly sample new configurations from a mixture of three Gaussian models, where each of the three Gaussian models are parameterized to be centered around the three best known configurations. To determine which configurations are best, we simply used the largest available fidelity budget evaluation for each of the configurations (see Algorithm 2).

We chose to use a mixture of three Gaussian models instead of a single Gaussian model, in case any of the low-fidelity budgets yield high-variance performance estimates that could cause the procedure to select the wrong configuration as the incumbent. We heuristically set the covariance of each of the Gaussian models to be a diagonal matrix parameterized such that the standard deviation of each hyper-parameter is equal to 5% of the total range of values for the hyper-parameter. Finally, each time a configuration needs to be suggested we sample a total of nine candidate configurations from the mixture of Gaussian models and suggest the one predicted to have the best performance by the surrogate model. The number of Gaussians, their standard deviations, and the number of samples drawn were each chosen based on some preliminary experiments on hand-crafted benchmarks.

For spline surrogate models, this procedure has the added benefit that we do not need to be able to locate the optimal configuration according the spline model, which may be non-trivial to do, as the model is not guaranteed to be uni-modal.

B Benchmark Scenarios

The first three scenarios we studied were also used in the original evaluation of BOHB (Falkner et al., 2018). The first two are for a Bayesian neural network trained using stochastic gradient Hamiltonian Monte-Carlo sampling (Chen et al., 2014) with scale adaption (Springenberg et al.,

Algorithm 2 The suggest function for our experimental AutoML configurators.

```

1: input
2:    $R$ , A dictionary mapping fidelities to configuration evaluation results
3:    $model$ , A dictionary mapping fidelities to previously-fit surrogate models (or None)
4:    $C$ , the parameter configuration space
5: output
6:    $c_{next}$ , a configuration to evaluate
7: procedure suggest
8:   # Find the largest-fidelity model, if any
9:   Initialize  $model_{current} := None$ 
10:  for each fidelity  $f \in R$  in descending order, do
11:    if  $model[f]$  is not None, then
12:       $model_{current} := model[f]$ 
13:    break
14:  # Sample a new configuration using the model, if possible
15:  if  $model_{current}$  is None, then
16:    Pick  $c_{next}$  by sampling a random configuration from  $C$ 
17:  else
18:    # Find the best configuration at any fidelity level
19:    Initialize an empty dictionary,  $performance := \{\}$ 
20:    for each fidelity  $f \in R$ , in ascending order, do
21:      for each configuration  $c \in R[f]$ , do
22:         $performance[c] := R[f][c]$ 
23:    # Sample from a mixture of 3 Gaussian models (see text for details)
24:    Sample 9 configurations from around the best 3 configurations in  $performance$ 
25:    Pick  $c_{next}$  as the best of the 9 configurations according to  $model_{current}$ 
26:  return  $c_{next}$ 

```

Table 2: The five real-world and four hand-crafted scenarios used to compare the four different hyper-parameter configurators.

Model/Algorithm	Dataset/Task	Fidelity Type	Fidelity Range	# HP
Bayesian Neural Network	Boston Housing	MCMC Steps	[500, 10 000]	7
	Protein Structure	MCMC Steps	[500, 10 000]	7
Proximal Policy Optimization	Cart Pole Swing-Up	# Training Runs	[1, 9]	5
Xgboost	King-Rook vs King-Pawn	# Estimators	[2, 128]	6
Histogram Gradient Boosting	Coverttype	Dataset Fraction	[1%, 100%]	7
Simulated Binary Classifier	Symmetric	Dataset Size	[500, 5 000]	1
	Asymmetric	Dataset Size	[500, 5 000]	1
	No Interactions	Dataset Size	[500, 5 000]	2
	Interactions	Dataset Size	[500, 5 000]	2

2016) on two UCI datasets (Dua and Graff, 2017), Boston housing and protein structure. We were unable to reproduce qualitatively similar results compared to those reported in Falkner et al. (2018) for either of these two scenarios. In each case, we observed that the gap between the performance for BOHB and Hyperband was much smaller than reported in their original study. From personal communication with the authors, we confirmed that we correctly set up the experiments and reproduced their method of analysis. To the best of our knowledge, the only differences in our execution environments were the particular machines used and, perhaps, the versions of some of the software packages.

The third scenario we used from their study configures the hyper-parameters of proximal policy optimization (Schulman et al., 2017) on the cart pole swing-up reinforcement learning task. In this case, even though the final performance gap between BOHB and Hyperband was also slightly smaller than originally reported, the results were still qualitatively similar.

Table 3: Results from applying the hyper-parameter configurators to the real-world AutoML scenarios. At each configuration budget we show the median test loss over the independent configurator runs with 95% bootstrap percentile confidence intervals. The median losses are shown in boldface if they are not worse than the best loss for a given configuration budget, according to a one-sided Welch t-test with a 5% significance level. All losses are scaled by dividing by the loss of the single best known configuration for each scenario. Smaller is better.

	BNN-BH	BNN-PS	PPO-CP	XGB-KR	HGB-CT
Budget (10%)	1 580 Seconds	2 420 Seconds	4 667 Seconds	148 Seconds	888 Seconds
CQA	1.91 [1.46, 2.42]	2.14 [1.63, 4.30]	10.49 [7.67, 13.74]	1.72 [1.44, 2.28]	1.28 [1.18, 1.35]
Spline	1.52 [1.40, 1.79]	1.77 [1.63, 1.87]	13.74 [9.42, 13.74]	1.72 [1.41, 2.41]	1.30 [1.21, 1.42]
Random (HB)	3.10 [1.60, 4.00]	2.13 [1.77, 4.74]	7.73 [6.21, 10.13]	1.91 [1.50, 2.00]	1.22 [1.13, 1.42]
TPE (BOHB)	1.76 [1.45, 2.05]	1.87 [1.56, 3.09]	7.70 [4.57, 9.84]	1.69 [1.47, 2.00]	1.23 [1.14, 1.48]
Budget (50%)	7 899 Seconds	12 102 Seconds	23 337 Seconds	742 Seconds	4 442 Seconds
CQA	1.44 [1.31, 1.67]	1.97 [1.72, 2.60]	5.73 [3.06, 6.47]	1.91 [1.44, 2.12]	1.17 [1.12, 1.24]
Spline	1.50 [1.42, 1.70]	1.85 [1.55, 2.24]	5.31 [4.10, 8.43]	1.56 [1.41, 1.94]	1.11 [1.10, 1.14]
Random (HB)	1.39 [1.36, 1.57]	1.50 [1.31, 1.83]	4.97 [3.25, 6.86]	1.78 [1.53, 1.97]	1.08 [1.07, 1.10]
TPE (BOHB)	1.38 [1.27, 1.68]	1.24 [1.10, 1.40]	3.59 [2.31, 5.37]	1.37 [1.28, 1.53]	1.09 [1.07, 1.14]
Budget (100%)	15 798 Seconds	24 204 Seconds	46 674 Seconds	1 484 Seconds	8 884 Seconds
CQA	1.44 [1.31, 1.67]	1.97 [1.62, 2.32]	4.45 [2.92, 6.11]	1.72 [1.41, 2.00]	1.12 [1.12, 1.17]
Spline	1.50 [1.43, 1.77]	1.85 [1.35, 2.24]	4.44 [3.53, 6.51]	1.56 [1.41, 1.94]	1.09 [1.06, 1.13]
Random (HB)	1.39 [1.36, 1.57]	1.50 [1.31, 1.83]	4.67 [2.69, 6.66]	1.72 [1.53, 1.97]	1.08 [1.06, 1.09]
TPE (BOHB)	1.38 [1.27, 1.68]	1.25 [1.10, 1.43]	2.22 [1.76, 3.21]	1.38 [1.31, 1.59]	1.05 [1.03, 1.08]

For the next two scenarios, we configured the hyper-parameters of Xgboost (Chen and Guestrin, 2016) on the UCI (Dua and Graff, 2017) king-rook vs king-pawn (Shapiro, 1987) dataset and histogram gradient boosting (scikit-learn’s (Pedregosa et al., 2011) implementation of a machine learning method inspired by LightGBM (Ke et al., 2017)) on the covtype dataset (Blackard and Dean, 1999).

For the simulated classifier scenarios, the classifier was constructed such that the variability in the measurements of the landscape mimic the behaviour of a binary classifier with a given error rate on a validation set of a given size. In particular, for each scenario, we defined functions that describe the landscape by mapping configurations to error rates, p , such that $0 < p < 1$. Then, given the error rate, p , of a particular configuration and a validation dataset size of n (the level of fidelity), we simulated an evaluation of a configuration by randomly drawing n trials from a binomial distribution with a success rate of p , and we recorded the loss as the number of successful trials divided by the total number of trials, n . We simulated short training times for the procedure by sleeping for one second per 1 000 instances in the dataset. This simulation ignores the variability in performance due to the random seed used when training a machine learning model; however, it still provides a cheap-to-evaluate and approximately-realistic method for generating landscapes with exactly known structure that can be used to evaluate the configurators.

C Extended Results

In Table 3, we show the results from applying the four hyper-parameter configuration procedures to the real-world AutoML scenarios. As can be seen, neither of the two new experimental methods provide results competitive with the two baselines: BOHB (Falkner et al., 2018) and Hyperband (Li et al., 2017). A similar analysis of other configuration budgets (not shown) yields qualitatively similar results.