A More details on DE

Differential Evolution (DE) is a simple, well-performing evolutionary algorithm to solve a variety of optimization problems [K. Price and Lampinen, 2006] [Das et al., 2016]. This algorithm was originally introduced in 1995 by Storn and Price [Storn and Price, 1997], and later attracted the attention of many researchers to propose new improved state-of-the-art algorithms [Chakraborty, 2008]. DE is based on four steps: initialization, mutation, crossover and selection. Algorithm 1 presents the DE pseudo-code.

Initialization. DE is a population-based meta-heuristic algorithm which consists of a population of \( N \) individuals. Each individual is considered a solution and expressed as a vector of \( D \)-dimensional decision variables as follows:

\[
\text{pop}_g = (x^1_i, x^2_i, ..., x^D_i), \quad i = 1, 2, ..., N, \quad (1)
\]

where \( g \) is the generation number, \( D \) is the dimension of the problem being solved and \( N \) is the population size. The algorithm starts initially with randomly distributed individuals within the search space. The function value for the problem being solved is then computed for each individual, \( f(x) \).

Mutation. A new child/offspring is generated using the mutation operation for each individual in the population by a so called mutation strategy. Figure 1 illustrates this operation for a 2-dimensional case. The classical DE uses the mutation operator \( \text{rand}/1 \), in which three random individuals/parents denoted as \( x_{r1}, x_{r2}, x_{r3} \) are chosen to generate a new vector \( v_i \) as follows:

\[
v_{i,g} = x_{r1,g} + F \cdot (x_{r2,g} - x_{r3,g}), \quad (2)
\]

where \( v_{i,g} \) is the mutant vector generated for each individual \( x_{i,g} \) in the population. \( F \) is the scaling factor that usually takes values within the range \( (0, 1] \) and \( r_1, r_2, r_3 \) are the indices of different randomly-selected individuals. Eq.2 allows some parameters to be outside the search range, therefore, each parameter in \( v_{i,g} \) is checked and reset\(^1\) if it happens to be outside the boundaries.

Crossover. When the mutation phase is completed, the crossover operation is applied to each target vector \( x_{i,g} \) and its corresponding mutant vector \( v_{i,g} \) to generate a trial vector \( u_{i,g} \). Classical DE uses the following uniform (binomial) crossover:

\[
\begin{aligned}
u_{i,g}^j &= v_{i,g}^j, \quad \text{if } (\text{rand} \leq p) \quad \text{or } (j = j_{\text{rand}}) \\
u_{i,g}^j &= x_{i,g}^j, \quad \text{otherwise}
\end{aligned} \quad (3)
\]

The crossover rate \( p \) is real-valued and is usually specified in the range \([0, 1]\). This variable controls the portion of parameter values that are copied from the mutant vector. The \( j \)th parameter value is copied from the mutant vector \( v_{i,g} \) to the corresponding position in the trial vector \( u_{i,g} \) if a random number is less than or equal to \( p \). If the condition is not satisfied, then the \( j \)th position is copied from the target vector \( x_{i,g} \). \( j_{\text{rand}} \) is a random integer in the range \([1, D]\) to ensure that at least one dimension is copied from the mutant, in case the random number generated for all dimensions is \( \geq p \). Figure 2 shows an illustration of the crossover operations.

Selection. After the final offspring is generated, the selection operation takes place to determine whether the target (the parent, \( x_{i,g} \)) or the trial (the offspring, \( u_{i,g} \)) vector survives to the next generation by comparing the function values. The offspring replaces its parents if it has a better\(^2\) function value.

---

\(^1\)A random value from \([0, 1]\) is chosen uniformly in this work

\(^2\)DE is a minimizer
on pre-determined geometrically spaced resource allocation. For DEHB we replace the random sampling with DE search. However, DEHB uses HB at its core to solve the “n versus B/n” tradeoff that HB was designed to address. Algorithm 2 shows how DEHB interfaces HB to query the sequence of how many configurations of each budget to run at each iteration. This view treats the DEHB algorithm as a sequence of predetermined (by HB), repeating Successive Halving brackets where, iteration number refers to the index of SH brackets run by DEHB.

Algorithm 2: A SH bracket under Hyperband

**Input:**
- \(b_{\text{min}}, b_{\text{max}}\) - min and max budgets
- \(\eta\) - fraction of configurations promoted
- iteration - iteration number

**Output:** List of no. of configurations and budgets

1. \(s_{\text{max}} = \lceil \log_{\eta} \frac{b_{\text{max}}}{b_{\text{min}}} \rceil\)
2. \(s = s_{\text{max}} - (\text{iteration mod } (s_{\text{max}} + 1))\)
3. \(N = \lceil \frac{s_{\text{max}} + 1}{\eta} \rceil\)
4. \(b_0 = \frac{b_{\text{max}}}{b_{\text{min}}} \cdot \eta^{-s}\)
5. budgets = n_configs = []
6. for \(i \in \{0, \ldots, s\}\) do
7. \(N_i = \lfloor N \cdot \eta^{-i} \rfloor\)
8. \(b = b_0 \cdot \eta^i\)
9. n_configs.append(N_i)
10. budgets.append(b)
11. end
12. return n_configs, budgets

C More details on DEHB

C.1 DEHB algorithm

Algorithm 3 gives the pseudo code describing DEHB. DEHB takes as input the parameters for HB \((b_{\text{min}}, b_{\text{max}}, \eta)\) and the parameters for DE \((F, p)\). For the experiments in this paper, the termination condition was chosen as the total number of DEHB brackets to run. However, in our implementation it can also be specified as the total absolute number of function evaluations, or a cumulative wallclock time as budget. L6 is the call to Algorithm 2 which gives a list of budgets which represent the sequence of increasing budgets to be used for that SH bracket. The nomenclature DE[budgets[i]], used in L9 and L12, indicates the DE subpopulation associated with the budgets[i] fidelity level. The if...else block from L11-15 differentiates the first DEHB bracket from the rest. During the first DEHB bracket (\(\text{iteration} == 0\)) and its second SH bracket onwards (\(\text{iteration} > 0\)), the top configurations from the lower fidelity are promoted\(^3\) for evaluation in the next higher fidelity. The function DE_trial_generation on L14, i.e., the sequence of mutation-crossover operations, generates a candidate configuration (config) to be evaluated for

\(^3\)only evaluate on higher budget and not evolve using mutation-crossover-selection
all other scenarios. L17 carries out the DE selection procedure by comparing the fitness score of config and the selected target for that DE evolution step. The target \((x_i, y_i\) from Equation 3) is selected on L9 by a rolling pointer over the subpopulation list. That is, for every iteration (every increment of \(j\)) a pointer moves forward by one index position in the subpopulation selecting an individual to be a target. When this pointer reaches the maximal index, it resets to point back to the starting index of the subpopulation. L18 compares the score of the last evaluated config with the best found score so far. If the new config has a better fitness score, the best found score is updated and the new config is marked as the incumbent, configinc. This stores the best found configuration as an anytime best performing configuration.

Algorithm 3: DEHB

\[
\begin{align*}
\text{Input:} & \quad b_{\min}, b_{\max} - \text{min and max budgets} \\
& \quad \eta - (\text{default } \eta=3) \\
& \quad F - \text{scaling factor (default } F = 0.5) \\
& \quad p - \text{crossover rate (default } p = 0.5) \\
\text{Output:} & \quad \text{Best found configuration, } configinc \\
1 & s_{\max} = \lfloor \log_\eta \frac{b_{\max}}{b_{\min}} \rfloor \\
2 & \text{Initialize } (s_{\max} + 1) \text{ DE subpopulations randomly} \\
3 & \text{bracket_counter} = 0 \\
4 & \text{while } \text{termination condition} \text{ do} \\
5 & \quad \text{for } \text{iteration} \in \{0, 1, ..., s_{\max} \} \text{ do} \\
6 & \quad \quad \text{budgets, } n_{\text{configs}} = \\
7 & \quad \quad \text{SH_bracket_under_HB}(b_{\min}, b_{\max}, \eta, \text{iteration}) \\
8 & \quad \quad \text{for } i \in \{0, 1, ..., s_{\max} - \text{iterations} \} \text{ do} \\
9 & \quad \quad \quad \text{target} = \text{rolling pointer for} \\
10 & \quad \quad \quad \text{DE[budgets[i]]} \\
11 & \quad \quad \quad \text{mutation_types} = \text{“vanilla” if } i \text{ is 0} \\
12 & \quad \quad \quad \text{else “altered”} \\
13 & \quad \quad \quad \text{if } \text{bracket_counter} \text{ is 0 and } i > 0 \text{ then} \\
14 & \quad \quad \quad \quad config = j\text{-th best config from } \\
15 & \quad \quad \quad \quad \text{DE[budgets[i - 1]]} \\
16 & \quad \quad \quad \text{else} \\
17 & \quad \quad \quad config = \text{DE_trial_generation(target,} \\
18 & \quad \quad \quad \text{mutation_type)} \\
19 & \quad \quad \text{result} = \text{Evaluate } config \text{ on} \\
20 & \quad \quad \text{budgets[i]} \\
21 & \quad \quad \text{DE selection using result, } config \text{ vs.} \\
22 & \quad \quad \text{target} \\
23 & \quad \quad \text{Update incumbent, } configinc \\
24 & \quad \text{end} \\
25 & \quad \text{bracket_counter} += 1 \\
26 & \text{end} \\
27 & \text{return } configinc
\end{align*}
\]

C.2 Handling Mixed Data Types

When dealing with discrete or categorical search spaces, such as the NAS problem, the best way to apply DE with such parameters is to keep the population continuous and perform mutation and crossover normally (Eq. 2, 3); then, to evaluate a configuration we evaluate a copy of it in the original discrete space as we explain below. If we instead dealt with a discrete population, then the diversity of population would drop dramatically, leading to many individuals having the same parameter values; the resulting population would then have many duplicates, lowering the diversity of the difference distribution and making it hard for DE to explore effectively. We designed DEHB to scale all parameters of a configuration in a population to a unit hypercube \([0, 1]\), for the two broad types of parameters normally encountered:

- **Integer and float** parameters, \(X^i \in [a_i, b_i]\) are retrieved as: \(a_i + (b_i - a_i) \cdot U_{1,9}\), where the integer parameters are additionally rounded.
- **Ordinal and categorical** parameters, \(X^i \in \{x_1, ..., x_n\}\), are treated equivalently s.t. the range \([0, 1]\) is divided uniformly into \(n\) bins.

We also experimented with another encoding design where each category in each of the categorical variables are represented as a continuous variables \([0, 1]\) and the variable with the max over the continuous variables is chosen as the category [Vallati et al., 2015]. For example, in Figure 3, the effective dimensionality of the search space will be 96-dimensional — 32 continuous variables + 64 continuous variables derived from 32 binary variables. We choose a representative set of benchmarks (NAS-Bench-201 and Counting Ones) to compare DEHB with the two encodings mentioned, in Figures 3, 4. It is enough to see one example which performs much worse than the DE-NAS [Awad et al., 2020] encoding we chose for DEHB. The encoding from [Vallati et al., 2015] did not achieve a better final performance than DEHB in any of our experiments.

C.3 Parallel Implementation

The DEHB algorithm is a sequence of DEHB Brackets, which in turn are a fixed sequence of SH brackets. This feature, along with the asynchronous nature of DE allows a parallel execution of DEHB. We dub the main process as the
**DEHB Orchestrator** which maintains a single copy of all DE subpopulations. An **HB bracket manager** determines which budget to run from which SH bracket. Based on this input from the bracket manager, the orchestrator can fetch a configuration\(^4\) from the current subpopulations and make an asynchronous call for its evaluation on the assigned budget. The rest of the orchestrator continues synchronously to check for free workers, and query the HB bracket manager for the next budget and SH bracket. Once a worker finishes computation, the orchestrator collects the result, performs DE selection and updates the relevant subpopulation accordingly. This form of an update is referred to as **immediate, asynchronous DE**.

DEHB uses a **synchronous** SH routine. Though each of the function evaluations at a particular budget can be distributed, a higher budget needs to wait on all the lower budget evaluations to be finished. A higher budget evaluation can begin only once the lower budget evaluations are over and the top \(1/\eta\) can be selected. However, the **asynchronous** nature of DE allows a new bracket to begin if a worker is available while existing SH brackets have pending jobs or are waiting for results. The new bracket can continue using the current state of DE subpopulations maintained by the **DEHB Orchestrator**. Once the pending jobs from previous brackets are over, the DE selection updates the **DEHB Orchestrator**’s subpopulations. Thus, the utilisation of available computational resources is maximized while the central copy of subpopulations maintained by the Orchestrator ensures that each new SH bracket spawned works with the latest updated subpopulation.

### D More details on Experiments

#### D.1 Baseline Algorithms

In all our experiments we keep the configuration of all the algorithms the same. These settings are well-performing setting that have been benchmarked in previous works — [Falkner et al., 2018], [Ying et al., 2019], [Awad et al., 2020].

**Random Search (RS)** We sample random architectures in the configuration space from a uniform distribution in each generation.

**BOHB** We used the implementation from https://github.com/automl/HpBandSter. In [Ying et al., 2019], they identified the settings of key hyperparameters as: \(\eta = 3\), the minimum bandwidth for the kernel density estimator is set to 0.3 and bandwidth factor is set to 3. In our experiments, we deploy the same settings.

**Hyperband (HB)** We used the implementation from: https://github.com/automl/HpBandSter. We set \(\eta = 3\) and this parameter is not free to change since there is no other different budget included in the NAS benchmarks.

**Tree-structured Parzen estimator (TPE)** We used the open-source implementation from https://github.com/hyperopt/hyperopt. We kept the settings of hyperparameters to their default.

**Sequential Model-based Algorithm Configuration (SMAC)** We used the implementation from https://github.com/automl/SMAC3 under its default parameter setting. Only for the Counting Ones problem with 64-dimensions, the **initial design** had to be changed to a Latin Hypercube design, instead of a Sobol design.

**Regularized Evolution (RE)** We used the implementation from [Real et al., 2019]. We initially sample an edge or operator uniformly at random, then we perform the mutation. After reaching the population size, RE kills the oldest member at each iteration. As recommended by [Ying et al., 2019], the population size (PS) and sample size (TS) are set to 100 and 10 respectively.

**Differential Evolution (DE)** We used the implementation from [Awad et al., 2020], keeping the **rand1 strategy for mutation** and **binomial crossover** as the crossover strategy. We also use the same population size of 20 as [Awad et al., 2020].

All plots for all baselines were plotted for the incumbent validation regret over the estimated wallclock time, ignoring the optimization time. The x-axis therefore accounts for only the cumulative cost incurred by function evaluations for each algorithm. All algorithms were run for similar actual wallclock time budget. Certain algorithms under certain benchmarks may not appear to have equivalent total estimated wallclock time. That is an artefact of ignoring optimization time. Model-based algorithms such as SMAC, BOHB, TPE have a computational cost dependent on the observation history. They might undertake lesser number of function evaluations for the same actual wallclock time.

#### D.2 Artificial Toy Function: Stochastic Counting Ones

![Figure 5: Results for the Stochastic Counting Ones problem for \(N = \{4, 8, 16, 32\}\) respectively indicating \(N\) continuous and \(N\) categorical hyperparameters for each case. All algorithms shown were run for 50 runs.](image)

The Counting Ones benchmark was designed to minimize the following objective function:

\[
f(x) = - \left( \sum_{x_i \in X_{cont}} x_i + \sum_{x_j \in X_{cont}} \mathbb{E}_b[(B_{y=x_j})] \right),
\]

where the sum of the categorical variables \((x_i \in \{0, 1\})\) represents the standard discrete counting ones problem. The continuous variables \((x_j \in [0, 1])\) represent the stochastic component with the budget \(b\) controlling the noise. The budget
here represents the number of samples used to estimate the mean of the Bernoulli distribution (B) with parameters \(x_j\).

The experiments on the Stochastic Counting Ones benchmark used \(N = \{4, 8, 16, 32\}\), all of which are shown in Figure 5. For the low dimensional cases, BOHB and SMAC’s models are able to give them an early advantage. For this toy benchmark the global optima is located at the corner of a unit hypercube. Random samples can span the lower dimensional space adequately for a model to improve the search rapidly. DEHB on the other hand may require a few extra function evaluations to reach similar convergence. However, this conservative approach aids DEHB for the high-dimensional cases where it is able to converge much more rapidly in comparison to other algorithms. Especially where SMAC and BOHB’s convergence worsens significantly. DEHB thus showcases its robust performance even when the dimensionality of the problem increases exponentially.

D.3 Feed-forward networks on OpenML datasets

![Figure 6: Results for the OpenML surrogate benchmark for the 6 datasets: Adult, Higgs, Letter, MNIST, Optdigits, Poker. The search space had 6 continuous hyperparameters. All plots shown were averaged over 50 runs of each algorithm.](image)

Figure 6, show the results on all 6 datasets from OpenML surrogates benchmark — Adult, Letter, Higgs, MNIST, Optdigits, Poker. The surrogate model space is just 6-dimensional, allowing BOHB and TPE to build more confident models and be well-performing algorithms in this space, especially early in the optimization. However, DE and DEHB are able to remain competitive and consistently achieve an improved final performance than TPE and BOHB respectively. While even TPE achieves a better final performance than BOHB. Overall, DEHB is a competitive anytime performer for this benchmark with the most robust final performances.

D.4 Bayesian Neural Networks

The search space for the two-layer fully-connected Bayesian Neural Network is defined by 5 hyperparameters which are: the step length, the length of the burn-in period, the number of units in each layer, and the decay parameter of the momentum variable. In Figure 7, we show the results for the tuning of Bayesian Neural Networks on both the Boston Housing and Protein Structure datasets for the 6-dimensional Bayesian Neural Networks benchmark. We observe that SMAC, TPE and BOHB are able to build models and reach similar regions of performance with high confidence. DEHB is slower to match in such a low-dimensional noisy space. However, given the same cumulative budget, DEHB achieves a competitive final score.

![Figure 7: Results for tuning 5 hyperparameters of a Bayesian Neural Network on the the Boston Housing and Protein Structure datasets respectively, for 50 runs of each algorithm.](image)

D.5 Reinforcement Learning

For this benchmark, the proximal policy optimization (PPO) [Schulman et al., 2017] implementation is parameterized with 7 hyperparameters: # units layer 1, # units layer 2, batch size, learning rate, discount, likelihood ratio clipping and entropy regularization. Figure 8 summarises the performance of all algorithms on the RL problem for the Cartpole benchmark. SMAC uses a SOBOL grid as its initial design and both its benefit and drawback can be seen as SMAC rapidly improves, stalls, and then improves again once model-based search begins. However, BOHB and DEHB both remain competitive and BOHB, DEHB, SMAC emerge as the top-3 for this benchmark, achieving similar final scores. We notice that the DE trace stands out as worse than RS and will explain the reason behind this. Given the late improvement for DE pop = 20, we posit that this is a result of the deferred updates of DE based on the classical DE [Awad et al., 2020] update design and also the design of the benchmark.

For classical-DE, the updates are deferred, that is the results of the selection process are incorporated into the population for consideration in the next evolution step, only after all the individuals of the population have undergone evolution. In terms of computation, the wall-clock time for population size number of function evaluations are accumulated, before the population is updated. In Figure 8 we illustrate why given how this benchmark is designed, this minor detail for DE slows down convergence. Along with a DE of population size 20 as used in the experiments, we compare a DE of population size 10 in Figure 8. For the Reinforcement Learning benchmark from [Falkner et al., 2018], each full budget function evaluation consists of 9 trials of a maximum of 3000 episodes. With a population of 20, DE will not
inject a new individual into a population unless all 20 individuals have participated as a parent in the crossover operation. This accumulates wallclock time equivalent to 20 individuals times 9 trials times time taken for a maximum of 3000 episodes. Which can explain the flat trajectories in the optimization trace for DE \( \text{pop} = 20 \) in Figure 8 (right). DE \( \text{pop} = 10 \) slashes this accumulated wallclock time in half and is able to inject newer configurations into the population faster and is able to search faster. Given enough runtime, we expect DE \( \text{pop} = 20 \) to converge to similar final scores. DEHB uses the immediate update design for DE, wherein it updates the population immediately after a DE selection, and not wait for the entire population to evolve. We posit that this feature, along with lower fidelity search, and performing grid search over population sizes with Hyperband, enables DEHB to be more practical than classical-DE.

![Figure 8](image.png)

Figure 8: (left) Results for tuning PPO on OpenAI Gym cartpole environment with 7 hyperparameters. Each algorithm shown was run for 50 runs. (right) Same experiment to compare DE with a population size of 10 and 20.

### D.6 NAS benchmarks

**NAS-Bench-101**

This benchmark was the first NAS benchmark relying on tabular lookup that was introduced to encourage research and reproducibility [Ying et al., 2019]. Each architecture from the search space is represented as a stack of cells. Each cell is treated as a directed acyclic graph (DAG) and the nodes and edges of these DAGs are parameterized which serve as the hyperparameters specifying a neural network architecture. NAS-Bench-101 offers a large search space of nearly 423k unique architectures that are trained on Cifar-10. The benchmark also offers a fidelity level — training epoch length — which allows HB, BOHB, and DEHB, to run on this benchmark. We run experiments on all 3 variants provided by NAS-101: Cifar A, Cifar B, Cifar C. The primary search space discussed by [Ying et al., 2019] is Cifar A; Cifar B and Cifar C are variants of the same search space with alternative encodings that deal with the hyperparameters defined on the edges of the DAG as categorical or continuous.

In the NAS-Bench-101 benchmark, the correlation between the performance scores and the different budgets are small [Ying et al., 2019], and therefore BOHB and DEHB do not yield better performance than the methods using full function evaluations only. All 3 evolutionary algorithms tested are able to exploit the discrete high-dimensional space much better than model-based methods such as BOHB and TPE, as seen by the performances of DE, DEHB and RE. While DEHB appears to be the algorithm with the best anytime performance in the high-dimensional discrete NAS space. DE yields the final best performance, closely followed by DEHB.

![Figure 9](image.png)

Figure 9: Results for Cifar A, B and C from NAS-Bench-101 for 26, 14, 27-dimensional spaces respectively. All algorithms reported for 50 runs.

**NAS-Bench-1shot1**

NAS-Bench-1shot1 was introduced by [Zela et al., 2020], as a benchmark derived from the large space of architectures offered by NAS-Bench-101. This benchmark allows the use of modern one-shot NAS methods with weight sharing ([Pham et al., 2018], [Liu et al., 2018]). The search space in NAS-Bench-1shot1 was modified to accommodate one-shot methods by keeping the macro network-level topology of the architectures similar and offering a different encoding design for the cell-level topology. This resulted in three search spaces: search space 1, search space 2 and search space 3 with 6240, 29160, and 363648 architectures respectively. In Figure 10, we show the results on all 3 search spaces. We exclude weight sharing methods from the algorithms compared, in order to maintain parity across all experiments, while focusing on the objective of comparing black-box solvers.

![Figure 10](image.png)

Figure 10: Results for the 3 search spaces from NAS-Bench-1shot1 for 50 runs of each algorithm. The 3 search spaces contains 9, 9, 11 categorical parameters respectively.

From among RS, TPE, SMAC, RE and DE — the full budget algorithms — only DE is able to improve significantly as training a single large architecture that contains all possible architectures in the search space.
optimization proceeds. For the multi-fidelity algorithms — HB, BOHB and DEHB — only DEHB is able to improve and diverge away from HB by the end of optimization. BOHB, HB, RS, TPE and RE all appear to follow a similar trace showing the difficulty of finding good architectures in this benchmark. Nevertheless, the DE-based family of algorithms is able to further exploit the search space and show better performance than the other algorithms. Though DE performs the best, RE remains competitive, again suggesting the power of evolutionary methods on discrete spaces. Among model-based methods, only TPE competes with DEHB.

**NAS-Bench-201**

To alleviate issues of direct applicability of weight sharing algorithms to NAS-Bench-101, [Dong and Yang, 2020] proposed NAS-Bench-201. This benchmark contains a fixed cell search space having DAGs with 4 nodes as the cell structure, and the edges of the DAG cells representing operations. The search space by design contains 6 discrete/categorical hyperparameters. NAS-Bench-201 provides a lookup table for 3 datasets: Cifar-10, Cifar-100 and ImageNet16-120, along with a fidelity level as number of training epochs. The search space for all 3 datasets include 15,625 cells/architectures. From the validation regret performances in Figure 11, it is clear that DEHB quickly converges to strong solutions which are a few orders of magnitude better than BOHB and RS (in terms of regret). DE and RE are both competitive with RE converging slightly faster than DE. Notably, DEHB is the only multi-fidelity algorithm in this experiment that works well.

NAS-Bench-201 specifies the same 6-dimensional discrete hyperparameter space for the Cifar-10 and Cifar-100 datasets. Figure 11 again shows that the evolutionary algorithms perform the best in a space defined by categorical parameters. SMAC in this scenario is the best-of-the-rest, outside of DEHB, RE and DE. BOHB evidently struggles to be even significantly better than HB.

**NAS-HPO-Bench**

To facilitate HPO research involving feed-forward neural networks, [Klein and Hutter, 2019] introduced NAS-HPO-Bench with a search space composed of hyperparameters that parameterize the architecture of a 2-layer feed-forward network\(^6\), along with hyperparameters for its training procedure. The primary difference between NAS-HPO from the OpenML surrogates benchmark is that in the latter, a random forest model was used as a surrogate to approximate the performance for configurations. NAS-HPO-Bench is designed in the same vein as the other NAS benchmarks discussed earlier. For the total of 9 discrete hyperparameters (4 for architecture + 5 for training), all 62208 configurations resulting from a grid search over the search space were evaluated to yield a tabular representation for configuration and performance mapping. The benchmark provides such lookup tables for 4 popular UCI regression datasets: Protein Structure, Slice Localization, Naval Propulsion and Parkinsons Telemonitoring. NAS-HPO-Bench also provides the number of training epochs as a fidelity level.

![Figure 11: Results for Cifar-10, Cifar-100, ImageNet16-120 from NAS-Bench-201 for 50 runs of each algorithm. The search space contains 6 categorical parameters.](image)

![Figure 12: Results for the Protein Structure, Slice Localization, Naval Propulsion, Parkinsons Telemonitoring datasets from NAS-HPO-Bench for 50 runs of each algorithm. The search space contains 9 hyperparameters.](image)

**D.7 Comparison of DEHB to BO-based multi-fidelity methods**

BOHB [Falkner et al., 2018] showed that its KDE based BO outperformed other GP-based BO methods. Hence, BOHB was treated as the primary challenger to DEHB as a robust, general multi-fidelity based HPO solver. In this section we run experiments on the benchmarks detailed in the previous sections, to compare DEHB to another popular multi-fidelity BO optimizer, Dragonfly [Kandasamy et al., 2020] (in addition to BOHB). Dragonfly implements BOCA (Kandasamy et al. [2017]) which performs BO with low-cost approximations of function evaluations on fidelities treated as a continuous

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\(^6\) additionally, a linear output layer
domain. However, this GP-based BO method had longer execution time compared to other algorithms for the tabular/surrogate benchmarks. In Figure 13 we therefore show average of 32 runs for each algorithm, while having to terminate runs earlier than other algorithms for certain cases. In this experiment, we optimize the median performance of a configuration over different seeds. We observe that Dragonfly shows a high variance in performance across benchmarks whereas DEHB is consistently the best or at worse, comparable to Dragonfly. Moreover, BOHB performs clearly better than Dragonfly in 8 out of the 16 cases shown in Figure 13, while being comparable to Dragonfly in at least 4 other benchmarks. Dragonfly comes out as the best optimizer only for the Cifar10 dataset in the NAS-201 benchmark in Figure 13. These experiments however, further illustrate the practicality, robustness, and generality of DEHB compared to GP-based multi-fidelity BO methods.

D.8 Results summary

In the previous experiments sections, results on all the benchmarks for DEHB and all other baselines were reported demonstrating the competitive and often superior anytime performance of DEHB. In Table 1, we report the mean final validation regret achieved by all algorithms across all the 26 benchmarks. DEHB got the best performance in nearly 1/3rd of the benchmarks while reporting the second-best performance in over 1/4-th of all the benchmarks. The last row of Table 1 shows the rank of each algorithm averaged across their final performances on each benchmark. DEHB clearly is the best performing algorithm on the whole, followed by DE, which powers DEHB under the hood. Such rankings illustrate DEHB’s robustness across different search spaces, including high dimensions, discrete or mixed type spaces, and even problems where response signals from lower fidelity subspaces may not be too informative. It must be noted that for all the different problems tested for with the collection of benchmarks, DEHB is never consistently outperformed by any multi-fidelity or full-budget algorithm.

It must be noted that based on the average rank plot in Figure 13, BOHB appears to be better than DEHB in the early middle section of the optimization. The underlying model-based search in BOHB can possibly explain this phenomenon. Though DEHB’s underlying DE requires more function evaluations to explore the space, it remains competitive with BOHB and the latter is not significantly better across any of the used benchmarks. Moreover, as Figure 13 indicates, BOHB’s relative performance worsens while DEHB continues to remain better even in comparison to the other full-budget black box optimizers such as DE, RE and TPE. BOHB’s model-based search can again be attributed for this phenomenon. Many of the benchmarks used are high-dimensional and have mixed data types, which can affect BOHB’s model certainty over the configuration space and require much more observations than DEHB requires. Overall, DEHB shows consistently good anytime performance with strong final performance scores too. As detailed earlier, DEHB’s efficiency, simplicity and its speed allow the good use of available resources and make it a good practical and reliable tool for HPO in practice.

Figure 13: Comparison of GP-based multi-fidelity BO (Dragonfly), KDE based multi-fidelity BO (BOHB), DE based multi-fidelity (DEHB) methods for some of the benchmarks, averaged over 32 runs of each algorithm.
Table 1: Final mean validation regret ± standard deviation for 50 runs all algorithms tested for DEHB is accompanied with its (rank) among other algorithms. The last row shows the average relative rank of each algorithm based on their final performance on each benchmark.

**E Ablation Studies**

DEHB was designed as an easy-to-use tool for HPO and NAS. This necessitated that DEHB contains as few hyperparameters as possible that require tuning or that which makes DEHB sensitive to them. Given that the HB parameters inherent to DEHB are contingent on the problem being solved, that leaves only the DE components’ hyperparameters to be set adequately. We perform ablation of the mutation and crossover rates to observe how it favors DEHB’s design for the suite of benchmarks we experiment on.

E.1 Varying F

The crossover probability $p$ was fixed at 0.5, while mutation factor $F$ was varied with the values 0.1, 0.3, 0.5, 0.7, 0.9. The studies are carried out on NAS-Bench-101, OpenML Surrogate and the toy Stochastic Counting Ones benchmarks. The results are reported in Figure 14.

We observe that a low $F$ of 0.1 allows more exploitative power to the DE search for a well correlated benchmark such as Counting Ones, while $F = 0.9$ performs the worst. However, for the other benchmarks $F = 0.1$ performs the worst with all other $F$ performing similarly. As a result we choose the conservative option of $F = 0.5$ to ensure one general set-
ting performs acceptably across all benchmarks.

E.2 Varying CR
The scaling factor $F$ was fixed at 0.5, while crossover factor $p$ was varied with the values 0.1, 0.3, 0.5, 0.7, 0.9. The studies were carried out on NAS-Bench-101, OpenML Surrogate and the toy Stochastic Counting Ones benchmarks. The results are reported in Figure 15.

The lower the $p$ value, the less likely are the random mutant traits to be incorporated into the population. For Counting Ones, we observe that a high $p$ slows down convergence, whereas low $p$ speeds up convergence. However, for the others, $p = 0.5$ is consistently the best performer. Hence, we chose $p = 0.5$ for the design of DEHB.
Figure 14: Ablation study for mutation factor $F$ for DEHB, with crossover probability fixed at $p = 0.5$
Figure 15: Ablation study for crossover probability $p$ for DEHB, with mutation factor fixed at $F = 0.5$
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