Can Subnetwork Structure be the Key to Out-of-Distribution Generalization?

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Abstract

Can models with particular structure avoid being biased towards spurious correlation in out-of-distribution (OOD) generalization? Peters et al. (2016) provides a positive answer for linear cases. In this paper, we use a functional modular probing method to analyze deep model structures under OOD setting. We demonstrate that even in biased models (which focus on spurious correlation) there still exist unbiased functional subnetworks. Furthermore, we articulate and demonstrate the functional lottery ticket hypothesis: full network contains a subnetwork that can achieve better OOD performance. We then propose Modular Risk Minimization to solve the subnetwork selection problem. Our algorithm learns the sub-network structure from a given dataset, and can be combined with any other OOD regularization methods. Experiments on various OOD generalization tasks corroborate the effectiveness of our method.

1. Introduction

Despite the remarkable progress we have witnessed in neural-network-based machine learning, the stories of failures continue to accumulate (Geirhos et al., 2020). Many of these failures are attributed to models exploiting spurious correlations or shortcuts (i.e. factors that are not used to generate the label). A colloquial example comes from (Beery et al., 2018) where the authors show how a neural network trained to distinguish cows from camels exploits shortcut such as background color for prediction. In a much more concerning example, (DeGrave et al., 2020) show how machine learning systems trained to detect COVID-19 exploited the data source (e.g., hospital) to artificially boost inference performance.

What causes these failures? Recent works (Peters et al., 2016; Arjovsky et al., 2019) argue the principle of empirical risk minimization (ERM) is at fault: if the data is generated from a fully observed causal bayesian network (CBN), then ERM would typically use all the features in the Markov blanket including those which are not the causes of the label. It may consequently fail to perform well under distribution shifts. This is known as the out-of-distribution (OOD) generalization problem. In an effort to alleviate the problem, Peters et al. (2016) proposes to first identify the target’s causal parents, and constrain the model structure by only updating the parameters for the parents. Nevertheless, their approach is only applicable for linear problem. Sagawa et al. (2020) also analyze the problem from a model structure perspective, but focusing on deep neural networks and showing that overparameterization will hurt OOD performance through data memorization and overfitting. Rather than focusing model structure, most recent works (Arjovsky et al., 2019; Sagawa et al., 2019; Ahuja et al., 2020; Krueger et al., 2020; Jin et al., 2020; Koyama & Yamaguchi, 2020; Creager et al., 2020) mainly target improvements in the objective function over ERM.

In this work, we set out to study the effect of the model structure in OOD generalization beyond simple considerations of model capacity. We begin by demonstrating that even already trained models that exploit spurious correlation can contain subnetworks that capture invariant features. We then
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We show that large trained networks that exploit spurious correlations contain subnetworks that are less susceptible to these spurious shortcuts.

We propose a novel functional lottery ticket hypothesis: there exists a subnetwork that can achieve better OOD and commensurate in-distribution accuracy in a comparable number of iterations when trained in isolation.

We propose Modular Risk Minimization (MRM), a straightforward and effective algorithm to improve OOD generalization. MRM helps select subnetworks and can be used in conjunction with other methods (e.g., IRM) and boosts their performance as well.

2. Invariant Prediction

2.1. Out-of-distribution (OOD) generalization problem

Consider a supervised learning setting where the data is gathered from different environments and each environment represents a different probability distribution. Let \((X^e, Y^e) \sim P^e\), where \(X^e \in \mathcal{X}, Y^e \in \mathcal{Y}\) stands for the feature random variable and the corresponding label, \(e \in \mathcal{E} = \{1, ..., E\}\) is the index for environments, and the set \(\mathcal{E}\) corresponds to all possible environments. The set \(\mathcal{E}\) is divided into two sets: seen environments \(\mathcal{E}_{\text{seen}}\) and unseen ones \(\mathcal{E}_{\text{unseen}} = \mathcal{E}_{\text{seen}} \cup \mathcal{E}_{\text{unseen}}\). The training dataset comprises samples from \(\mathcal{E}_{\text{seen}}\). The dataset from environment \(e\) is given as \(D^e = \{(x^e_i, y^e_i)\}_{i=1}^{n^e}\), where each point \((x^e_i, y^e_i)\) is an independently identically distributed (IID) sample from \(P^e\) and \(n^e\) is the number of samples in environment \(e\). We write the training dataset as \(D_{\text{train}} = \bigcup_{e \in \mathcal{E}_{\text{seen}}} D^e\). In the rest of the work, we interchangeably use the term domain and environment, and we will use in-distribution or in-domain to refer to seen environmental data, and out-distribution or out-domain for unseen environmental data.

Let \(f_\theta : \mathcal{X} \rightarrow \mathcal{Y}\) denote the parametrized model with parameters \(\theta \in \Theta\). Define the risk achieved by the model as \(\mathcal{R}^e(\theta) = \mathbb{E}^e[\ell(X^e, Y^e)]\) where \(\ell\) is the loss per sample (e.g., cross-entropy, square loss). The goal of out-of-distribution (OOD) generalization problem is to learn a model that solves

\[
\min_{\theta \in \Theta} \max_{e \in \mathcal{E}} \mathcal{R}^e(\theta).
\]

Since we only have access to data from \(D_{\text{train}}\) and do not see samples from the unseen environments, the above problem can be challenging to solve.

**Data generation process.** We assume \(X^e\) is generated from latent variables \(Z^e = (Z^e_{\text{inv}}, Z^e_{\text{sp}})\). Consider an illustrative example where \(X^e\) could be the pixels in images, while \(Z^e_{\text{inv}}\) denotes invariant features (e.g., foreground) and \(Z^e_{\text{sp}}\) denotes spurious features (e.g., background). We write \(X^e = G(Z^e_{\text{inv}}, Z^e_{\text{sp}})\), where \(G\) is a map from the latent space to the pixel space. \(Y^e\) is the label for the object and it is determined based on the following map \(Y^e = F(Z^e_{\text{inv}})\). The combination pattern of \(Z^e_{\text{inv}}\) and \(Z^e_{\text{sp}}\) varies across domains, hence generating different environmental distributions.

In our description of the data generation, we do not use noise variables to keep things simple (\(Y\) is related to \(Z\) deterministically and \(X\) is related to \(G\) deterministically). Suppose that we can recover \(Z^e_{\text{inv}}\) and \(Z^e_{\text{sp}}\) from \(X^e\) and we write these inverse maps as \(Z^e_{\text{inv}} = G^*_{\text{inv}}(X^e)\) and \(Z^e_{\text{sp}} = G^*_{\text{sp}}(X^e)\). The ideal function that the model wants to learn is \(F \circ G^*_{\text{inv}}\) as it yields zero error and only relies on invariant latents.

However, as we explain next that due to selection biases the model can often find it hard to learn a model that only relies on \(Z^e_{\text{inv}}\).

**Bias.** To explain why the datasets have a bias, let us consider a simple example, where \(Z^e_{\text{inv}} \in \{-1, 1\}^{D_{\text{inv}}}\), \(Z^e_{\text{sp}} \in \{-1, 1\}^{D_{\text{sp}}}\) and \(Y^e \in \{-1, 1\}\). Suppose each component of \(Z^e_{\text{inv}}\) is \(Y^e\) and each component of \(Z^e_{\text{sp}}\) independently takes a value equal to \(Y^e\) with a probability \(p^e\) and \(-Y^e\) with a probability \(1 - p^e\). If \(p^e\) is close to 1 and \(G^*_{\text{sp}}\) is an easier function to learn than \(G^*_{\text{inv}}\), then it is intuitive that the model can instead learn \(Z^e_{\text{sp}}\) and predict the label \(Y^e\). However, this can be catastrophic as the correlation between the spurious feature and the label only holds in the training environments and does not translate to the test environments where \(p^e = \frac{1}{2}\).

Even if \(p^e\) is small, as long as \(Z^e_{\text{sp}}\) is high dimensional \((D_{\text{sp}} >> D_{\text{inv}})\), the model can be shown to significantly rely on \(Z^e_{\text{sp}}\) (Nagarajan et al., 2020). The above example uses binary valued latents for ease of exposition, but the same biases can occur in more general settings where the same problems plague the models.

2.2. A Motivating Example

In this section, we use a simple example to motivate the constraints we impose in our approach. Consider the data setting described in the previous section, \(Z^e_{\text{inv}} \in \{-1, 1\}\) \((D_{\text{inv}} = 1)\) and \(Z^e_{\text{sp}} \in \{-1, 1\}^{D_{\text{sp}}}\) \((D_{\text{sp}} = D)\). We take \(G\) to be
the identity map as in Tsipras et al. (2019); Rosenfeld et al. (2020) and thus $X^e = (Z^e_{inv}, Z^e_{sp})$. Suppose the model $f_\theta$ is a linear predictor; we refer to the components associated with invariant features as $w_{inv}$ and those associated with the spurious feature as $w_{sp}$.

**Learning a sparse classifier:** Find a maximum margin classifier that satisfies the following sparsity constraint: the number of non-zero coefficients $\leq d$. We denote such a classifier as $f^d_{sparse}$.

In the next proposition, we compare the behavior of the sparse classifier that we defined above with a classifier that relies only on spurious features. We construct a regular classifier $f_{reg}$ (with unit norm) that purely relies on the spurious features, i.e., $w_{inv} = 0$ and $w_{sp} = 1/\sqrt{D_{sp}}$ and thus has poor OOD performance. We denote the average error rate of the classifier $h$ on seen (or unseen) environments as $Err_{seen}(h)$ (or $Err_{unseen}(h)$). Here the error for binary classification is defined to be $Err(h) = \frac{1}{2}E_{(X', Y') \sim p}[1 - Y' h(X')]$. We denote the margin of classifier for data in environment $e$ as Margin$^e$.

**Proposition 1.** Consider the dataset in Section 2.1 with $Z^e_{inv} \in \{−1, 1\} (D_{inv} = 1)$ and $Z^e_{sp} \in \{−1, 1\}^D (D = D_{sp})$. Let $n$ be the number of training samples in $D_{train}$, $c$ be a constant in $(0, 1)$ such that for all $e \in E_{seen}$, $p^e > \frac{1}{2} + \frac{c}{2}$ and $p^e < \frac{1}{2}$ for $e \in E_{unseen}$. For sparsity constraint $d = 2$, we have:

- **Compare margin for in-distribution sample:** for any $\delta \in (0, 1)$, if $D \geq \frac{1}{c^2} \sqrt{2ln(n)/\delta}$, then with a probability at least $1 - \delta$, Margin$^e_{seen}(f^d_{sparse}) < \text{Margin}^e_{seen}(f_{reg})$.

- **Similar in-distribution performance:** $\forall e \in E_{seen}$, $Err^e_{seen}(f^d_{sparse}) = 0$, $Err^e_{seen}(f_{reg}) \leq 2e^{-2c^2D}$.

- **Better out-distribution performance:** $\forall e \in E_{unseen}$, $Err^e_{unseen}(f^d_{sparse}) = 0$ and $Err^e_{unseen}(f_{reg}) = 0.5$.

From the above Proposition, we can conclude that if $c$ or $D$ is high, then the train accuracy of the sparse classifier and the regular classifier are similar but the OOD accuracy of the two classifiers are different with sparse classifier being much better. The algorithm is likely to select the regular classifier over the sparse classifier as it has a much higher margin than the sparse classifier.

Proposition 1 compares the optimal sparse classifier with a purely spurious one. Both have same in-distribution performance, but the former has a better OOD performance. We compare the margins to show that if we use a gradient descent on logistic loss, it will be biased towards the spurious classifier (Soudry et al., 2018). We clarify that Proposition 1 is not intended to show a tradeoff between OOD performance and margin. Consider the experiment of spiral vs. linear boundary of Sec 3.1 in Parascandolo et al. (2020). In the experiment, the spiral boundary is associated with invariant features and the linear boundary is associated with spurious ones. The authors set the margin for linear boundary to be larger than the that of the spiral boundary. In this case, ERM learns a model that uses spurious features. Even if we were to reduce the margin of the linear boundary to be smaller than the spiral boundary, ERM continues to rely on the spurious features as it prefers to use a simpler margin (Shah et al., 2020).

For this linear setting that we discussed above, we can learn a constrained max-margin classifier by adding $\ell_1$ constraints. This is a tractable problem to solve as the problem remains convex. However, as we move to neural networks, learning sparse classifiers with good OOD performance is significantly more challenging owing to the non-convexity. This issue is the subject of later sections. Before we address this issue of learning sparse networks with good OOD properties, there is another important question to be answered. In the setting of the above proposition, we rely on the fact that a sparse model exists that relies on invariant features only and yields better OOD performance. How do we do know that this is a proper assumption for real datasets used for neural network training? In the next section, we analyze neural networks via modular subnetwork introspection to show that such a sparse model exists.

### 3. A Functional Modularity Based Analysis

#### 3.1. Preliminaries

**Technical approach.** The modularity property of neural network has long been considered as an essential foundation of systematical generalization (Ballard, 1987; Marcus, 1998; Csordás et al., 2020). Consider a task that can be compositionally separated into different independent subtasks, we aim to probe a functional module subpart of the full neural network that can solve one particular subtask. Following Zhou et al. (2019); Csordás et al. (2020), we identify different subnetworks which perform different functions, from a given pretrained network.

Specifically, we deem functional modules to be particular subsets of the weights inside a neural network. For a $L$ layer neural network model $f(w_1, \cdots, w_L; \cdot)$ where $\theta = \{w_1, \cdots, w_L\}$, we model the subnetwork with a set of binary masks $m_1 \in \{0, 1\}^{n_1}$ on the $l$-th layer weight tensor $w_l \in \mathbb{R}^{n_l}$, where $n_l$ is the number of dimensionality of the $l$-th layer network parameters. The subnetwork is then given by $f(m_1 \odot w_1, \cdots, m_L \odot w_L; \cdot)$. Further, in order to make this subnetwork structure learnable, we assume each entry of the mask to be independent Bernoulli random variables, and model their logits as $\pi_l \in \mathbb{R}^{n_l}$. Hence, in this probabilistic modeling setting, the $l$-th layer subnetwork structure $m_l$ is generated by performing Bernoulli
We use the term F modularity probing and thereby uncover the corresponding functional modules. We will interchangeably use the term of module within the neural network in the form of binary tensor $\{m_1, \cdots, m_L\}$, and thereby uncover the corresponding functional module within the neural network in the form of binary tensor $m = \{m_1, \cdots, m_L\} = \{\text{sigmoid}(\pi_l) > 0.5 | l = 1, 2, \cdots\}$. Using the term modularity probing method to refer to this technique subsequently. We will interchangeably use the term of module and subnetwork due to their consistency in our context.

Dataset construction. We take the intuition from Arjovsky et al. (2019); Nam et al. (2020); Ahuja et al. (2021); Ahmed et al. (2021) to design a biased variant of the original MNIST dataset (LeCun et al., 1998). A discussion about the difference between ours and theirs is deferred to supplementary materials. The digit shape semantics are considered as $Z_{\text{inv}}$ while color semantics as $Z_{\text{sp}}$. We choose ten different kinds of color and define a one-to-one corresponding bias relationship with ten digit class (e.g., “2” ↔ “green”, “4” ↔ “yellow”). For each domain, we define the bias coefficient to be the ratio of the data that obeys this relationship. Those images which don’t follow this relationship are then assigned with random colors. The bias coefficient for two in-domains is (1.0, 0.9) respectively, which means the first domain is completely biased and 90 percent of the second domain is biased. For the out-domain, all images are assigned a random color for evaluating to how much extent the model has learned the invariant feature. The out-domain will serve as a tool environment for module learning in this section, representing a thorough disentanglement of two attributions. It will then act as the test distribution in a realistic setting in Section 5. Unless otherwise specified, the label is set as the class where the invariant attribution lies. We use the term FULLCOLOREDMNIST to refer to this task to distinguish with the binary colored mnist dataset in Arjovsky et al. (2019).

Algorithms analyzed. We study four OOD generalization algorithms in this paper: Empirical Risk Minimization (ERM) (Vapnik, 1999), Invariant Risk Minimization (IRM) (Arjovsky et al., 2019), Risk Extrapolation (REx) (Krueger et al., 2020) and group Distributional Robust Optimization (DRO) (Sagawa et al., 2019). More details about them are left to supplementary materials. Figure 2(a) plots the generalization performance of these algorithms w.r.t. the training process. REx (76.17%) and DRO (78.56%) methods surpass ERM baseline by a large margin, while IRM (59.55%) only gets slightly better results than ERM (58.04%). The failure of IRM in realistic problems has been analyzed in Jin et al. (2020); Nagarajan et al. (2020); Rosenfeld et al. (2020); Ahuja et al. (2021) and attributed to the overparameterization regime and curse of dimensionality, hence we omit related discussion here.

3.2. Modular subnetwork introspection

Departing from previous approaches, in this section we think of learning the digit and color semantics as different functional subtasks of the original task, rather than opposite non-spurious / spurious features. We split the out-domain into two parts and refer to them as the in-split and out-split of the out-domain (terminology from Gulrajani & Lopez-Paz (2020)). We define two subtasks, identification of digit and identification of color. For each subtask, we assume that we have access to respective semantic labels. It’s important to note that the semantic color label is used here for analysis and is not a part of our main method described later.

In order to study the functional module for the two subtasks, we apply the modularity probing method to diagnose given pretrained models. Specifically, we separately train and get a digit and a color subnetwork for each model across different algorithms and training steps. We evaluate the obtained digit modules’ behaviors on the out-split of out-domain (as the in-split has been taken for module searching). Figure 2(b) suggests a significant evidence that, even for biased models such as ERM trained ones, there exist unbiased invariant subnetworks (digit modules) with good OOD generalization ability. We also explore this property for other modules and algorithms and defer these results to supplementary materials.

Discussion about the sparsity of digit weights / features. We additionally visualize the Bernoulli probability of learned subtask modules. Figure 3 displays the first layer of model trained on two in-domains with ERM. We can see that the color feature is more pervasive than digit feature, spreading over a broader range across the neurons. Although the sparsity of weights is not exactly the sparsity of features, the discovery is aligned with the assumption in Proposition 1 that $D_{\text{inv}}$ has a small number of dimensionality. The visualization results of other layers are similar to
The visualization of the Bernoulli probability of digit and color functional module for the first (convolutional) layer. The weight tensor is reshaped to two dimension for display convenience. The probability takes value from 0 to 1.

In this section we confirm in large trained models, there lie invariant functional modules, viz. subnetwork that behaves well for target invariant function (e.g., digit classification). However, it’s more worthwhile to find out about whether an appropriate subnetwork structure can help improve in the sense of OOD generalization during training. We investigate this problem in the next section.

**Figure 3.**

**4. Structure Matters: Towards Functional Lottery Tickets Hypothesis**

Frankle & Carbin (2018) proposes the lottery ticket hypothesis from a pruning perspective, suggesting that among all different subnetworks, there exists a so-called “winning ticket” that can reach the generalization ability of the full network with faster training speed. In this original lottery ticket hypothesis, the data distribution remains unchanged across training and testing. Whereas in our OOD context, we seek a model subnetwork whose functional predictions are invariant with respect to the change in distribution.

**The functional lottery ticket hypothesis:** A randomly initialized, dense neural network contains a subnetwork that is initialized such that — when trained in isolation — it can achieve better out-of-distribution performance w.r.t. the given function (e.g., digit identification in our context) than the original full network after training for the same number of iterations.

Concretely, our functional lottery ticket hypothesis claims that for a dense neural network model \( f(w; x) \) with initialization parameter \( w^0 \), there exists a module \( m \) enabling a subnetwork \( f(m \odot w; x) \) to surpass the OOD performance of full network when trained from \( f(m \odot w^0; x) \) on in-distribution data. Note that this is a stronger statement than the original lottery ticket hypothesis, which only requires the winning ticket to reach similar performance to the full network in the IID setting.

**Demonstration.** How do we identify the functional winning tickets? How should one search for a structure that is best for OOD generalization? To unravel the possible best result one can reach, we design an “oracle” subnetwork. After obtaining an ERM trained model, the structure of oracle module is found with the aid of the in-split data part of out-domain. Namely, we use the modularity probing technique introduced in Section 3.1 with these “oracle” data from the out-distribution, and then deploy the resulting subnetwork back onto the previous initialization. We then evaluate all methods on the out-split.

According to Sagawa et al. (2020), the underparameterized regime can keep the model from overfitting to spurious features. Therefore, we choose the random subnetwork as another option to investigate whether sparsity / underparametrization alone can achieve an unbiased solution. “randwhole” method keeps the ratio of full network same as the oracle subnetwork. In other works it has been claimed that the sparsity per layer is the only working factor for pruning (Su et al., 2020; Frankle et al., 2020), we also experiment with the “randlayer” method, where we randomly sample subnetworks with the same per-layer-sparsity.

Figure 1 and 4(b) show the OOD and in-distribution generalization results for these subnetworks with ERM training on the FULLCOLOREDMNIST dataset respectively. The oracle subnetwork beats the original ERM by a large margin for OOD and maintains indistinguishable performance for the in-distribution examples, confirming that a good module structure can indeed surpass the full network in terms of this digit function. We show their performance with greater levels of sparsity in Figure 4(a) and see a considerable consistent accuracy gap between oracle and random baselines for all level of sparsity. The validity of our functional lottery ticket hypothesis is thereby empirically affirmed, and we thus propose that appropriate structure induction can impose a needed inductive bias to prevent the model from fitting the spurious correlation. We also conclude that sparsity constraint imposed cannot help alone, as two random
methods don’t yield non-trivial benefit than ERM (both under 60% accuracy). Additionally, we demonstrate our hypothesis is also applicable for other OOD algorithms in the supplementary materials.

Discussion. One can rightly criticize this investigation as unfair in that it compares a method using out-domain data to baselines without such privileged access. We acknowledge this issue and, for now, only seek out this “oracle subnet-work” to highlight the importance of structure. We now turn to the question of how we can design a practical structure searching algorithm to overcome this limitation.

4.1. Modular risk minimization

The motivation behind the Modular Risk Minimization (MRM) method is to get rid of spurious features by hunting for a desired functional winning ticket. Since we have shown in previous subsection that contrary to Sagawa et al. (2020), only sparsity constraints imposed at the beginning of training cannot do the magic, we propose the following criterion:

A good structure should balance the predictiveness for invariant feature and sparsity well.

Our procedure first trains the model with ERM resulting in a potentially biased classifier. At this time, the functional lottery ticket hypothesis suggests that the model has already learned a promising functional module within. Hence, we simply apply the subnetwork probing technique with training data to learn the potential advantageous structure. The structure learning objective takes a combination of cross entropy loss and sparsity regularization to balance the two desiderata mentioned above. We then simply train from scratch again only with the weights in the obtained subnetwork and fix the other weights to zero. We summarize this procedure in Algorithm 1, where i, c, l are respectively the index for a datum, label class and network layer. It’s notable that in Figure 1, our proposed MRM algorithm successfully reaches a very close accuracy to the oracle optimal structure.

Structure learning by invariance capturing. Notably, MRM does not impose any invariance across domains and is thus orthogonal to the advantage of other OOD algorithms. Unlike heuristic structure searching paradigms (Lee et al., 2018; Wang et al., 2020), our method can incorporate any OOD generalization approach to help improve the structure learning and model training process. This enables MRM to act as a plug-in method to boost other algorithms by supplying a good subnetwork learned from their respective objectives. We simply replace the cross entropy loss \( \mathcal{L}_{CE} \) in Algorithm 1 with recently developed OOD losses: \( \mathcal{L}_{IRM}, \mathcal{L}_{REx} \) and \( \mathcal{L}_{DRO} \) by IRM, REx and DRO. These new methods are therefore referred by Modular Invariant Risk Minimization (ModIRM), Modular Risk Extrapolation (ModREx) and Modular Distributionally Robust Optimization (ModDRO). With these explicitly OOD learning algorithms, the cross domain variance is taken into account and thereby ameliorates the invariance property of the subnetwork. We note that more flexible combinations can be explored (e.g., use a different loss design for subnetwork and model learning), but we leave this for future work and only study these three variants in this paper.

4.2. Ablation for winning tickets learning

To better understand the crucial succeeding reasons for two kinds of winning tickets – oracle subnetworks and MRM subnetworks, we conduct corresponding ablation studies.

Importance of initialization. One of the main argument of Frankle & Carbin (2018) is that the winning tickets cannot be learned effectively without its original initialization. We verify this for our hypothesis as well. Figure 5(a) depicts the failure of functional winning tickets when random re-initialization is performed before the training of subnetworks. At this time, both subnetworks achieve similar OOD performance to full network ERM. This ablation study confirms the importance of reusing initialization.

Effects of bias relationship. Our FullColoredMNIST keeps one fixed color-digit relationship for all biased data. Will rearranging this bias relationship defined in Section 3.1 before the subnetwork is trained destroy the tickets? We then

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**Algorithm 1 Modular Risk Minimization**

**Input:** Data \( \{(x_i^e, y_i^e)\}_{i,e} \), neural network \( f(w; \cdot) \), subnetwork logits \( \pi \), the coefficient of sparsity penalty \( \alpha \), number of steps for model and subnetwork structure training \( N_1, N_2 \).

**Stage 1:** full model (pre-) train

Get model initialization \( w_0 \).

for \( n=1 \) to \( N_1 \) do

\( \text{Update} \ f \text{ with } \mathcal{L}_{CE}(w) := \sum_{i,c}y_{i,c} \log f(w; x_i)_c. \)

end for

**Stage 2:** module structure probing

for \( n=1 \) to \( N_2 \) do

Sample subnetwork \( m \sim \text{sigmoid}(\pi) \).

Update module \( \pi \) with

\[ \mathcal{L}_{MOD} = \mathcal{L}_{CE}(m \odot w) + \alpha \sum_{l,j} \pi_{l,j}. \]

end for

**Stage 3:** subnetwork retrain

Obtain the module by hard thresholding:

\( m = \{ \pi_l > 0 \mid l = 1, 2, \ldots \} \).

Set model parameters back to \( w_0 \).

for \( n=1 \) to \( N_1 \) do

\( \text{Update} \ f \text{ with } \mathcal{L}_{CE}(m \odot w). \)

end for
We compare our algorithm and its OOD variants with recent methods aiming at robust predictions across environments. In this section, we demonstrate the effectiveness of our modular risk minimization algorithm on a variety of datasets. We keep the same object (invariant feature) to the images across different columns. For all methods we keep the same model architecture and training settings. We build three OOD datasets according to the bias protocol introduced in Section 2.1: FULL COLORED MNIST, COLORED OBJECT and SCENE OBJECT. For all datasets we design two training in-domains and one out-domain for evaluating OOD generalization capability. We defer all other experimental details to the supplementary materials.

FULL COLORED MNIST. Details of the construction are in Section 2.1. We summarize the results in Table 1. We also use ERM trained with completely unbiased data whose bias coefficient is (0,0,0,0,0) to serve as an upper bound (coined as “Unbias” in the tables). Our method can consistently promote the OOD performance on this task, bringing around 10% accuracy promotion. The best behaved algorithm, ModDRO, reaches 85.53% accuracy, contrary to the 78.25% of top-grade baseline DRO and 99.0% achieved by unbiased solution.

COLORED OBJECT. We take inspiration from Ahmed et al. (2021) to build this biased dataset together with the following SCENE OBJECT one. Ten classes of objects extracted from MSCOCO dataset (Lin et al., 2014) are put onto ten kinds of color backgrounds. Figure 6(a) displays 100 samples from this crafted biased dataset. Like FULL COLORED MNIST, we also set a one-to-one object-color relationship and set the bias coefficient differently as (0.8,0.6,0.0). Results in Table 2 demonstrate the advantages of our methods: all our four methods all achieve accuracy above 50%, boosting their different baselines towards the optimal “unbias” solution.

SCENE OBJECT. Ten classes of objects extracted from MSCOCO dataset are put onto ten kinds of scenery backgrounds from Places dataset (Zhou et al., 2018). These scenery backgrounds make this task a more complex one than COLORED OBJECT. Figure 6(b) displays 100 samples

| METHODS | TRAIN ACCURACY | TEST ACCURACY |
|---------|----------------|---------------|
| ERM     | 87.56 ± 2.52   | 43.74 ± 2.11  |
| MRM     | 94.01 ± 0.82   | 54.85 ± 2.11  |
| IRM     | 88.68 ± 2.11   | 45.4 ± 2.40   |
| ModIRM  | 93.01 ± 0.36   | 52.35 ± 1.28  |
| REX     | 89.85 ± 1.50   | 47.20 ± 3.43  |
| ModREX  | 93.55 ± 1.45   | 55.51 ± 2.76  |
| DRO     | 91.73 ± 0.40   | 51.95 ± 1.62  |
| ModDRO  | 92.67 ± 0.92   | 55.20 ± 1.40  |
| UNBIAS  | 95.00 ± 0.70   | 72.37 ± 2.53  |

| METHODS | TRAIN ACCURACY | TEST ACCURACY |
|---------|----------------|---------------|
| ERM     | 98.10 ± 0.09   | 57.75 ± 1.84  |
| MRM     | 98.90 ± 0.05   | 72.98 ± 0.58  |
| IRM     | 98.18 ± 0.09   | 59.30 ± 1.88  |
| ModIRM  | 98.77 ± 0.12   | 70.86 ± 2.12  |
| REX     | 98.86 ± 0.10   | 75.61 ± 1.26  |
| ModREX  | 99.28 ± 0.04   | 82.06 ± 0.73  |
| UNBIAS  | 99.07 ± 0.04   | 99.03 ± 0.08  |
from this crafted dataset. Like FULLCOLOREDMNIST, we set a one-to-one object-scenery relationship and set the bias coefficient to be \((0.9, 0.7, 0.0)\), making it a even more biased and thus more difficult one than the previous task. This can also be shown with only 56.46% accuracy of unbiased solution. Corresponding results in Table 3 shows that for this highly biased task, MRM and its variants can still accordingly improve out-distribution generalization performance in this highly bias setting, where previous OOD algorithms bring very limited benefit.

### 6. Related Work

**Out-of-distribution generalization.** Machine learning beyond IID assumption is a very important problem and many research areas such as domain adaptation (Crammer et al., 2008; Ben-David et al., 2010) and domain generalization (Muandet et al., 2013; Motiian et al., 2017) have received much attention (Gulrajani & Lopez-Paz, 2020). To get stable prediction for new unseen data distribution, it is desired to only rely on invariant features among the causal factorization of physical mechanisms of problem settings (Schöllkopf et al., 2012). Peters et al. (2016) (ICP) claims that the residual of invariant method should remain IID and thus proposes to adopt statistical tests for mining invariant feature set. Rojas-Carulla et al. (2018) generalizes this approach to nonlinear settings.

Recently, since Arjovsky et al. (2019) brings invariant prediction into a more practical scenario, a large amount of works has made solid progress for alleviating spurious correlation and shortcut exploitation (Geirhos et al., 2020; Koh et al., 2020); Sagawa et al. (2019) proposes to use group DRO when attribution information is provided; Chang et al. (2020) incorporates this invariant inference idea into selective rationalization area; Ahuja et al. (2020) studies the IRM formulation from a game theory and bilevel optimization formulation; Krueger et al. (2020) propose REX to enforce the variance of losses across distribution, which is further analyzed by Xie et al. (2020a); Koyama & Yamaguchi (2020) (IGA) also has a similar contribution with different theoretical analysis; Jin et al. (2020) (RGM) proposes another training objective from regret minimization viewpoint; Pezeshki et al. (2020) studies the gradient starvation phenomenon which is connected with spurious correlation and proposes an insightful solution; Creager et al. (2020) (EIL) points out that invariant prediction shares the same spirit with fair representation learning; Parascandolo et al. (2020) (ILC) proposes to focus second order landscape information; Ahmed et al. (2021) adopts a divergence term to match the output distribution spaces of different domains; Müller et al. (2020) achieves invariance from an information theory start point and enforces conditional invariance with HSIC terms. Some other works also point out the pitfalls of current approaches, showing only in very limited situations can Arjovsky et al. (2019) (e.g., low dimension settings) really capture invariance: Rosenfeld et al. (2020) proves the validity of IRM for linear cases but gives a negative example for nonlinear cases; Nagarajan et al. (2020) analyzes different failure modes of OOD generalization; Ahuja et al. (2021) analyze the sample efficiency properties of IRM; Kamath et al. (2021) investigates the success and failure cases of IRM and IRMv1 on simple but insightful settings, and claims the community might need a better invariance notion.

Another line of works study a related but different topic named debiasing, where there is no explicit multiple environments setting provided. Bias in realistic datasets are usually exploited in a spurious way, such as the texture-bias of Imagenet-trained models (Geirhos et al., 2018). Subsequent works (Wang et al., 2019; Bahng et al., 2020; Shi et al., 2020; Nam et al., 2020; Li et al., 2021; Sauer & Geiger, 2021) focus on addressing the bias problem with explicit debiasing procedure.

**Modularity.** Modularity (Ballard, 1987; Fodor et al., 1988; Newman, 2006) has been considered as a crucial part of intelligent systems. Lots of works focus on imposing explicit module level modularity (Clune et al., 2013; Andreas et al., 2016; Chang et al., 2018; Goyal et al., 2021), while others also explore weight level modularity in a more fine-grained way (Mallya & Lazebnik, 2018; Watanabe et al., 2019; Filan et al., 2020; Csordás et al., 2020). Our work also belongs to the latter category.

**Pruning.** We mainly focus on unstructured pruning literature. This line of model compression literature dates back to Mozer & Smolensky (1989); LeCun et al. (1989); Hassibi & Stork (1993) with more recent pruning methods (Han et al., 2015; Molchanov et al., 2016; Dong et al., 2017). Recently, the lottery ticket hypothesis (Frankle & Carbin, 2018) sheds more light into this field, showing the importance of initialization. (Liu et al., 2018) also propose another viewpoint that for practical settings the inherited weights are not important.

### Table 3. Generalization performance on SCENEOBJECT.

| METHODS | TRAIN ACCURACY | TEST ACCURACY |
|---------|----------------|---------------|
| ERM     | 98.87 ± 0.23   | 37.29 ± 2.74  |
| MRM     | 99.61 ± 0.04   | 39.44 ± 0.77  |
| IRM     | 98.68 ± 0.27   | 37.19 ± 2.58  |
| ModIRM  | 99.39 ± 0.01   | 39.14 ± 1.34  |
| REX     | 92.91 ± 1.11   | 38.84 ± 1.39  |
| ModREX  | 96.71 ± 0.53   | 41.04 ± 1.46  |
| DRO     | 98.89 ± 0.35   | 36.34 ± 1.67  |
| ModDRO  | 99.41 ± 0.13   | 39.14 ± 1.60  |
| UNBIAS  | 95.25 ± 2.21   | 56.46 ± 0.75  |
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7. Discussion

Data settings. The seminal work (Arjovsky et al., 2019) proposes to use color in digit identification as a spurious correlation. In order to exploit the effectiveness of IRM, the authors enforce a 25% label noise in the binary classification data and assign color a larger correlation than the true digit shape. In this way, ERM exploits color feature to predict. While there is controversy surrounding whether one should still treat digit as desired learning target under this situation, we choose to impose no label noise in FULL-COLOREDMNIST as is the case in Nam et al. (2020); Ahmed et al. (2021). This choice enables the structure learning procedure could mine the true invariant feature. On the other hand, our work is limited as we haven’t considered the data settings such as group attribution available ones (Sagawa et al., 2019; Xie et al., 2020b; Khani & Liang, 2021) and we shall fill this gap in future work. More about datasets can be found in Section C.1.

Success of MRM. There are several reasons for why MRM can improve OOD performance without invariance constrain. The first reason is related to our label noise free setting discussed above. This makes the invariant feature itself perfectly predictive of the label, thus containing all information about the desired target function. Then the problem would be how to exploit this information effectively. MRM becomes competent for OOD tasks by providing a novel and helpful parameterization method for the original optimization problem with extra parameters. One notable thing is that more structure parameters actually don’t increase the expressive power of the neural network, since every weight can take zero value in nature. Another reason is we adopt an explicit approach to zero out the “spurious part” of the model weights, hence achieving a not-so-biased solution. Notice this cannot be achieved with random sparse model, revealing the structure is a key element for OOD generalization. Therefore, a positive answer is given to the title of this work. We further refer to Section C.3 for empirical results of the importance of a proper sparsity level in structure learning.

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A. Omitted Proof

A.1. Proof for Proposition 1

Proof. We first analyze the performance of \( f_{\text{reg}} \). The prediction from this classifier is \( \hat{Y}^c = \text{sgn}(w_{\text{sp}}^T Z^c_{\text{sp}}) \), thus

\[
w_{\text{sp}}^T Z^c_{\text{sp}} = \frac{1}{\sqrt{D}} \sum_{i=1}^{D} Z^c_{\text{sp},i} = \sqrt{D} \left[ \frac{1}{D} \sum_{i=1}^{D} Z^c_{\text{sp},i} \right],
\]

\[
\hat{Y}^c = \text{sgn}(w_{\text{sp}}^T Z^c_{\text{sp}}) = \text{sgn}\left[ \frac{1}{D} \sum_{i=1}^{D} Z^c_{\text{sp},i} \right].
\]

(2)

We then analyze the error for any environment:

\[
\text{Err}^e = \frac{1}{2} \left[ 1 - \mathbb{E}^e[\hat{Y}^c Y^c] \right],
\]

\[
\mathbb{E}^e[\hat{Y}^c Y^c] = \mathbb{E}^e[\text{sgn}\left( \frac{1}{D} \sum_{i=1}^{D} Z^c_{\text{sp},i} \right) Y^c] = \sum_{y \in \{-1, 1\}} \mathbb{P}[Y^c = y] \mathbb{E}^e[\text{sgn}\left( \frac{1}{D} \sum_{i=1}^{D} Z^c_{\text{sp},i} \right) Y^c = y],
\]

(3)

\[
\mathbb{E}^e[\text{sgn}\left( \frac{1}{D} \sum_{i=1}^{D} Z^c_{\text{sp},i} \right) Y^c = 1] = \mathbb{P}\left[ \frac{1}{D} \sum_{i=1}^{D} Z^c_{\text{sp},i} > 0 | Y^c = 1 \right] - \mathbb{P}\left[ \frac{1}{D} \sum_{i=1}^{D} Z^c_{\text{sp},i} \leq 0 | Y^c = 1 \right]
\]

\[
= 2\mathbb{P}\left[ \frac{1}{D} \sum_{i=1}^{D} Z^c_{\text{sp},i} > 0 | Y^c = 1 \right] - 1.
\]

(4)

Observe that \( \mathbb{P}\left[ \frac{1}{D} \sum_{i=1}^{D} Z^c_{\text{sp},i} < 0 | Y^c = 1 \right] = \mathbb{P}\left[ \frac{1}{D} \sum_{i=1}^{D} Z^c_{\text{sp},i} > 0 | Y^c = -1 \right] \). Using this observation and plugging equation 4 into equation 3 we get

\[
\text{Err}^e = \mathbb{P}\left[ \frac{1}{D} \sum_{i=1}^{D} Z^c_{\text{sp},i} \leq 0 | Y^c = 1 \right].
\]

(5)

Let us now bound \( \text{Err}^e \). Define \( Z^c_{\text{sp}} = \frac{1}{D} \sum_{i=1}^{D} Z^c_{\text{sp},i} \). Since

\[
\mathbb{E}^e\left[ \frac{1}{D} \sum_{i=1}^{D} Z^c_{\text{sp},i} | Y^c = 1 \right] = 2p^c - 1,
\]

we have

\[
\mathbb{P}\left[ \frac{1}{D} \sum_{i=1}^{D} Z^c_{\text{sp},i} \leq 0 | Y^c = 1 \right] = \mathbb{P}\left[ \bar{Z}^c_{\text{sp}} \leq 0 | Y^c = 1 \right]
\]

\[
= \mathbb{P}\left[ \bar{Z}^c_{\text{sp}} - \mathbb{E}\left[ \bar{Z}^c_{\text{sp}} \right] \leq -\mathbb{E}\left[ \bar{Z}^c_{\text{sp}} \right] | Y^c = 1 \right] \leq \mathbb{P}\left[ \bar{Z}^c_{\text{sp}} - \mathbb{E}\left[ \bar{Z}^c_{\text{sp}} \right] \geq \mathbb{E}\left[ \bar{Z}^c_{\text{sp}} \right] | Y^c = 1 \right]
\]

\[
\leq 2e^{-2(2p^c - 1)^2 D} \leq 2e^{-2c^2 D}.
\]

(7)

In the test environment, since \( Z^c \) and \( Y^c \) are independent and \( p^c = 0.5 \). As a result, the error in test environment for the regular classifier is \( \text{Err}^e = \mathbb{P}\left[ \frac{1}{D} \sum_{i=1}^{D} Z^c_{\text{sp},i} \leq 0 | Y^c = 1 \right] = \mathbb{P}\left[ \frac{1}{D} \sum_{i=1}^{D} Z^c_{\text{sp},i} \leq 0 \right] = 0.5 \).

Now let us consider the optimal sparse max-margin classifier. For \( d = 2 \), the max-margin classifier for the above data distribution is simply \( w_{\text{inv}} = 1 \) and \( w_{\text{sp}} = 0 \). Since \( Z^c_{\text{sp}} = Y^c \), in both train and test, the sparse classifier has a perfect accuracy in both train and test environments.

Next, we compare the margins. The margin for \( f_{\text{reg}} \) is \( Y^c w_{\text{reg}} Z_{\text{sp}} \geq \frac{c}{\sqrt{D}} \) with a probability at least \( 1 - \delta \) (the proof follows directly from Hoeffding’s inequality and we refer to the Appendix A of Nagarajan et al. (2020)). In comparison, \( f^d_{\text{sp}} \) that assigns weights to invariant and spurious parts as follows \( w_{\text{inv}} = 1 \) and \( w_{\text{sp}} = 0 \) achieves a margin of \( Y^c w_{\text{inv}} Z_{\text{inv}} = 1 \).
B. More Discussion

One possible algorithm we do not explore in this work is the IRM games (Ahuja et al., 2020). This approach sees the IRM formulation from a game theory perspective across different environments and design a corresponding algorithm. The algorithm has one network per environment thus the number of parameters scale in number of environments. In order to make comparisons apple to apple, we keep all the methods to have the same parameter complexity. IRM games will have more parameters so we do not compare with it. Other methods including extra parameters (e.g., auxiliary neural networks) are also not considered due to analogous reason and we shall explore them in future work.

We notice there is a deep connection between Parascandolo et al. (2020) and our work. Our MRM explicitly learn a subnetwork architecture and only update the corresponding subset of weights, while Parascandolo et al. (2020) also restricts its optimization within a subset of weights by “and mask” algorithm, which only updates a parameter when gradients across domains are consensus to each other. However, this method may need a appropriate large number of domains to work (see details in their paper) and hence is not among the studied algorithms in this paper. In the future work, we intend to explore the relationship between the structure of our digit module and the subset found by their “and mask” under more domains.

Similar to our method, Dropout (Srivastava et al., 2014) also only updates part of the network during training. We list some of the main difference here: Dropout aims to prevent overfitting by simply not updating whole model parameters towards one single function, while we pursue to identify one particular functional module architecture within full network; Dropout only randomly zero out the updating gradients, while we intentionally pick particular functional part of the model in an end-to-end way; what’s more, Dropout is not activated during testing inference time, while our subnetwork is kept for all following stages.

In Hooker et al. (2019; 2020), the authors propose that model compression will hurt the accuracy on underrepresented groups with negligible impacts on overall accuracy, which seems contradictory to our results. Here we state about the difference in the settings and claim that their works are actually consistent in spirit to ours. First of all, their work do not target OOD problems or invariant prediction, but focus on long tail underrepresented subgroups in IID situations. In our context “bias” means a spurious / shortcut way of inference and we aim to zero out the spurious part in the model parameters, while their settings don’t contain a spurious feature that will bias the model prediction, hence most of the parameters are rightful and shouldn’t be got rid of. As a result, it’s natural for model compression to hurt in their cases. What’s more, our method utilize a much more careful subnetwork selection method where we aim to maximize the in-domain performance when searching structures, ensuring that we do not wipe off the useful and rightful part of parameters.

In Bengio et al. (2019) and its following analysis (Priol et al., 2020), the authors claim that a better knowledge of true causal relationship will bring a faster speed of transferring learning. While on the other hand, from Figure 1 we can see a good subnetwork leads to faster convergence and better performance. Although we do not target a causally promising algorithm in this work (our main goal is to get rid of spurious correlation for OOD problems), we claim our approach is connected with causal discovery (Bengio et al., 2019) stated above, in the sense of capturing causal relationship hidden in data better (please note that invariant prediction can be seen as a higher level of transfer learning in an OOD sense).

C. More on Experiments

C.1. Dataset details

**FULLCOLOREDMNIST.** We use ten colors taken from Ahmed et al. (2021) for all the data. Their RGB values are: [0, 100, 0], [188, 143, 143], [255, 0, 0], [255, 215, 0], [0, 255, 0], [65, 105, 225], [0, 225, 225], [0, 0, 255], [255, 20, 147], [160, 160, 160]. Each image is of size $3 \times 32 \times 32$ and the whole dataset contains 60000 images. The difference between Ahmed et al. (2021) and ours are significant: their whole data is deemed to come from “majority group” and “minority group”, where the images from “majority group” are colored with previous mentioned ten colors and those from “minority group” are colored with other fifty different colors. In contrast, all of our images are colored with these ten colors. We set a bias relationship to connect each digit and each color one by one and define bias coefficient to be the ratio of bias data that follows the relationship, as described in Section 3.1. Besides, their data are pooled and integrated as one domain before presented to algorithms. Our dataset is also different with Arjovsky et al. (2019), where they use label noise intentionally to make the correlation between label and color to be higher than the one between label and digit. This makes ERM severely biased towards color and thus fails in test domain. On the contrary, we do not impose any sort of label noise and leave label as the ground truth digit information. What’s more, Arjovsky et al. (2019) use binary classification and two colors, which is a much simpler setting. Our setting is also different from Nam et al. (2020), which is not multi environment and only has
Table 4. Generalization performance on FULLCOLOREDMNIST with oracle validation.

| METHODS | TRAIN ACCURACY | TEST ACCURACY |
|---------|----------------|---------------|
| ERM     | 98.10 ± 0.10   | 58.04 ± 1.95  |
| MRM     | 98.90 ± 0.05   | 73.21 ± 0.58  |
| IRM     | 98.17 ± 0.12   | 59.55 ± 1.90  |
| MODIRM  | 98.67 ± 0.20   | 70.35 ± 3.22  |
| REX     | 98.83 ± 0.09   | 76.17 ± 1.53  |
| MODREX  | 99.28 ± 0.05   | 82.13 ± 0.82  |
| DRO     | 98.94 ± 0.11   | 78.56 ± 1.42  |
| MODDRO  | 99.38 ± 0.06   | 85.67 ± 0.51  |
| UNBIAS  | 99.05 ± 0.04   | 97.86 ± 0.20  |

Table 5. Generalization performance on COLOREDOBJECT with oracle validation.

| METHODS | TRAIN ACCURACY | TEST ACCURACY |
|---------|----------------|---------------|
| ERM     | 87.58 ± 2.42   | 44.39 ± 2.44  |
| MRM     | 94.00 ± 0.56   | 55.03 ± 1.76  |
| IRM     | 87.63 ± 2.32   | 44.49 ± 2.15  |
| MODIRM  | 92.95 ± 0.41   | 52.55 ± 0.65  |
| REX     | 89.28 ± 1.35   | 46.07 ± 2.59  |
| MODREX  | 82.65 ± 1.85   | 55.56 ± 3.16  |
| DRO     | 91.84 ± 2.17   | 53.20 ± 1.15  |
| MODDRO  | 92.32 ± 1.37   | 55.46 ± 1.43  |
| UNBIAS  | 92.32 ± 1.80   | 72.77 ± 3.48  |

one domain.

COLOREDOBJECT. We imitate Ahmed et al. (2021) to take ten objects as invariant features and put them on ten different color backgrounds. We use the ten colors mentioned above, and take the following ten objects: boat, airplane, truck, dog, zebra, horse, bird, train, bus, motorcycle. Each image is of size $3 \times 64 \times 64$ and the whole dataset contains 10000 images. The difference between ours and Ahmed et al. (2021) is similar to FULLCOLOREDMNIST: their settings contain a minority group which has many other backgrounds than the mentioned ten color backgrounds.

SCENEOBJECT. We imitate Ahmed et al. (2021) to take ten objects as invariant features and put them on ten different scenery backgrounds (beach, canyon, building facade, staircase, desert sand, crevasse, bamboo forest, broadleaf, ball pit and kasbah). We use the same ten object classes mentioned above. Each image is of size $3 \times 64 \times 64$ and the whole dataset contains 10000 images. The difference between ours and Ahmed et al. (2021) is similar to FULLCOLOREDMNIST: their settings contain a minority group which has many other backgrounds than the mentioned ten scenery backgrounds.

C.2. Experimental details

For the results in the main text, we take a commonly used policy to report the last step accuracy. We also apply the another evaluation method (the “oracle validation” in Gulrajani & Lopez-Paz (2020)) to report accuracy, and provide corresponding results in Table 4, 5 and 6. These results are consistent to those in main text, showing the validity of our conclusion. For both method, we take a similar approach to Gulrajani & Lopez-Paz (2020), and the difference is that we take a finite search set for hyperparameters instead of sampling of a human defined distribution. For all datasets, we search the regularization coefficient of IRM and REx in $\{1e^{-1}, 5e^{-1}, 1, 1e1, 1e2, 1e3, 1e4\}$, the step when the regularization is added into training in $\{0, 1000, 2000\}$. Furthermore, we also search a binary option about whether to scale down the whole loss term by the regularization coefficient as in Arjovsky et al. (2019). For DRO we search the group proportion step size $\eta_q$ (notation taken
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Table 6. Generalization performance on SCENEOBJECT with oracle validation.

| METHODS   | TRAIN ACCURACY | TEST ACCURACY |
|-----------|----------------|---------------|
| ERM       | 95.04 ± 2.13   | 37.44 ± 1.15  |
| MRM       | 98.83 ± 0.26   | 39.49 ± 0.15  |
| IRM       | 92.66 ± 0.24   | 37.54 ± 0.94  |
| ModIRM    | 94.63 ± 0.54   | 40.04 ± 1.90  |
| REX       | 92.73 ± 1.61   | 39.39 ± 0.96  |
| ModREx    | 96.72 ± 0.53   | 40.79 ± 1.62  |
| DRO       | 94.53 ± 2.17   | 36.64 ± 1.35  |
| ModDRO    | 93.52 ± 5.90   | 40.99 ± 1.70  |
| UNBIAS    | 85.47 ± 2.37   | 56.91 ± 1.31  |

from Sagawa et al. (2019)) in \{1e \to 4, 1e \to 3, 1e \to 2, 1e \to 1, 1\}. For the subnetwork structure learning, we follow Csordás et al. (2020) to use Adam optimizer and search the logit learning rate among \{1e \to 2, 1e \to 1, 1\} and the sparsity coefficient among \{1e \to 8, 1e \to 7, 1e \to 6, 1e \to 5, 1e \to 4, 1e \to 3\}. For Figure 4(a), the x-axis is in log scale and ranges from 1e \to 5 to 1e \to 3 for sparse cases.

For FULLCOLOREDMNIST, all experiments are measured by computing mean and standard deviation across five trials with random seeds. For optimization we use SGD + momentum (0.9) with 1e \to 4 weight decay, where the initial learning rate is 1e \to 1 and is decayed every 600 steps for all algorithms. We take the batch size to be 128. The training process longs for 2000 steps (and thus N1 is also 2000). N2 is set to 2000, and it can be seen from Table 7 that the importance of this hyperparameter is very limited. We use a simple ConvNet with three convolutional layers with feature map dimensions of 64, 128 and 256, each followed by a ReLU nonlinear and a batch normalization layer. The fourth layer is a fully connected layer. For COLOREDOBJECT, all experiments are measured by computing mean and standard deviation a cross three trials with random seeds. The learning rate is 1e \to 1 and decays every 1200 steps. The training process longs for 2000 steps. Others are kept the same with previous dataset. We use Wide ResNet 28-2 architecture (Zagoruyko & Komodakis, 2016). To adapt to 64 \times 64 image size, we replace the average pooling layer with window size 8 to one with size 16. For SCENEOBJECT, we use a learning rate of 5e \to 2 which decays every 1200 steps. Other unmentioned settings are kept consistent as above. We use Tesla V100 GPU to perform the experiments.

C.3. More empirical results

We affirm that color is more fitted to neural network prior by plotting the digit accuracy of digit module and color accuracy of color module w.r.t. the training process in Figure 7(a). The two modules are both learned given the same ERM trained model at every step. The result shows that the training of color module converges much faster than the digit one, and thus showing that neural networks have a natural favor for color (texture) information than digit (shape) information as shown in Zeiler & Fergus (2014); Ritter et al. (2017); Brendel & Bethge (2019); Shi et al. (2020). We also plot the behaviors of digit module learning of three OOD algorithms in Figure 7(b) 7(c) and 7(d). These three plots show a very similar results.
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Figure 8. The digit / color accuracy for some logical operation results of learned digit and color module.

to Figure 2(b). We omit the behavior of color module of these algorithms since it’s very similar to that of ERM shown in Figure 7(a).

We play with the obtained digit and color module with several basic logical operations in Figure 8. The intersection of digit and color module is important as its complement behaves trivially in Figure 8(e) and 8(f). However, the intersection module alone can not express predictiveness (Figure 8(a)) for digit identification. We show that it needs to be combined with other part to work in Figure 8(g) and 8(h). We also additionally visualize the two modules for the linear layer in Figure 9(a) and get similar results to that in Figure 3.

Figure 9. Compensatory experimental results.

For completeness we also compare with magnitude pruning, although it’s not for out-of-distribution generalization and its structure search process is data independent like us. The results for ERM on FULLCLOREDMNIST are in Figure 9(b). We can see our methods surpass this widely used pruning approach (e.g., by the original lottery ticket hypothesis paper and most prevailing approaches (Han et al., 2015; Zhu & Gupta, 2017)) in all level of sparsity. Please see more related discussion in Section C.4. Furthermore, we do ablation about the hyperparameters of subnetwork learning in the stage 2 of MRM algorithm in Table 7 to show the importance of proper level of “good” sparsity (instead of random sparsity, whose importance is restricted as discussed in main text). In Table 7 we also show results with different values of $N_2$ in MRM algorithm. Although $N_2$ we adopt is not the optimal setting, we simply use it since it’s enough to achieve an approximate
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Table 7. Ablation for hyperparameters of subnetwork learning. All numbers are out-distribution generalization accuracy.

| LR  | α   | 1e−3   | 1e−4   | 1e−5   | 1e−6   | 1e−7   | N2   | ACCURACY     |
|-----|-----|--------|--------|--------|--------|--------|------|-------------|
| 1e − 2 | 1e−3 | 50.45 ± 2.70 | 68.56 ± 0.39 | 67.46 ± 0.42 | 64.13 ± 1.05 | 63.25 ± 1.12 | 1000 | 71.46 ± 1.89 |
| 1e − 1 | 1e−3 | 36.95 ± 2.25 | 76.83 ± 0.81 | 72.98 ± 0.58 | 71.91 ± 0.85 | 71.30 ± 0.55 | 2000 | 72.98 ± 0.58 |
| 1   | 1e−3 | 21.76 ± 1.29 | 50.19 ± 2.33 | 65.54 ± 2.16 | 66.31 ± 2.56 | 65.51 ± 3.31 | 3000 | 73.39 ± 0.84 |
| 1e−1 | 1e−1 | 36.95 ± 2.25 | 76.83 ± 0.81 | 72.98 ± 0.58 | 71.91 ± 0.85 | 71.30 ± 0.55 | 5000 | 73.70 ± 0.59 |

Table 8. Ablation for the third stage of MRM.

| METHODS | w/ stage 3 | w/o stage 3 |
|---------|------------|-------------|
| MRM     | 72.98 ± 0.58 | 62.99 ± 1.96 |
| ModIRM  | 70.86 ± 2.12 | 58.87 ± 2.40 |
| ModREX  | 82.06 ± 0.73 | 77.48 ± 2.30 |
| ModDRO  | 85.53 ± 0.61 | 80.47 ± 1.87 |

convergence of stage 2. We further do ablation about MRM for with or without stage 3 in Table 8, showing the importance of the last stage and the validity of our functional lottery ticket hypothesis.

C.4. Towards a different pruning method

In this paper we propose an OOD algorithm coined as MRM. Here we point out that MRM can actually be seen as a different pruning method and can be applied to IID tasks as well. We do not state this explicitly in the main text since it’s not closely related to our OOD research category. We slightly modify MRM to serve as a pruning algorithm in this way: we also use three stages training paradigm as MRM in main text, but specifically for stage 2, we jointly train the subnetwork structure and model parameters, and let the algorithm break the looping once the specified sparsity level is reached. We shall not claim this method to be a novel pruning method as we notice that it’s similar (though stille different) to Louizos et al. (2018), which for unknown reasons is not taken into baseline consideration by recent pruning works. We now show related experimental results in the sense of pruning settings (which is IID generalization) in Table 9. The codes and baseline results are based on Wang et al. (2020). We also keep the hyperparameters setting to be consistent with Wang et al. (2020), and take the modular learning rate 0.01 and modular sparsity regularization 0.0001. Our method outperforms other baseline methods, especially for extremely sparse 98% pruning ratio cases. We think of this as an insightful advantage of our approach.

Table 9. Our pruning method on CIFAR10 with ResNet32 across different pruning ratios.

| METHODS      | RATIO | 90%  | 95%  | 98%  |
|--------------|-------|------|------|------|
| FULL NETWORK |       | 94.23| –    | –    |
| OBD (LeCun et al., 1989) | 94.17 | 93.29 | 90.32 |
| M. Prune (Zeng & Urtasun, 2018) | 94.21 | 93.02 | 90.31 |
| LT (Frankle & Carbin, 2018) | 92.31 | 91.06 | 88.78 |
| LT Rewind (Frankle et al., 2020) | 93.97 | 92.46 | 89.18 |
| DSR (Mostafa & Wang, 2019) | 92.97 | 91.61 | 88.46 |
| SET (Mocanu et al., 2018) | 92.30 | 90.76 | 88.29 |
| Deep-R (Mocanu et al., 2018) | 91.62 | 89.84 | 86.45 |
| SNIP (Lee et al., 2018) | 92.59 | 91.01 | 87.51 |
| GRASP (Wang et al., 2020) | 92.38 | 91.39 | 88.81 |
| OURS         |       | 94.19 | 93.36 | 92.80 |