Generalizing Goal-Conditioned Reinforcement Learning with Variational Causal Reasoning

Abstract

As a pivotal component to attaining generalizable solutions in human intelligence, reasoning provides great potential for reinforcement learning (RL) agents' generalization towards varied goals by summarizing part-to-whole arguments and discovering cause-and-effect relations. However, how to discover and represent causalities remains a huge gap that hinders the development of causal RL. In this paper, we augment Goal-Conditioned RL (GCRL) with *Causal Graph (CG)*, a structure built upon the relation between objects and events. We novelly formulate the GCRL problem into variational likelihood maximization with CG as latent variables. To optimize the derived objective, we propose a framework with theoretical performance guarantees that alternates between two steps: using interventional data to estimate the posterior of CG; using CG to learn generalizable models and interpretable policies. Due to the lack of public benchmarks that verify generalization capability under reasoning, we design nine tasks and then empirically show the effectiveness of the proposed method against five baselines on these tasks. Further theoretical analysis shows that our performance improvement is attributed to the virtuous cycle of causal discovery, transition modeling, and policy training, which aligns with the experimental evidence in extensive ablation studies.

1 Introduction

The generalizability, which enables an algorithm to handle unseen tasks, is fruitful yet challenging in multifarious decision-making domains. Recent literature [1, 2, 3] reveals the critical role of reasoning in improving the generalization of reinforcement learning (RL). However, most off-the-shelf RL algorithms [4] have not regarded reasoning as an indispensable accessory, thus usually suffer from data inefficiency and performance degradation due to the mismatch between training and testing settings. To attain generalization at the testing stage, some efforts were put into incorporating domain knowledge to learn structured information, including sub-task decomposition [5], or program generation [6, 7, 8] and execution [9, 10], which guide the model to solve complicated tasks in an explainable way. However, such symbolism-dominant methods heavily depend on the re-usability of sub-tasks and pre-defined grammars, which may not always be accessible in decision-making tasks.

Inspired by the close link between reasoning and the cause-and-effect relationship, causality is recently incorporated to compactly represent the aforementioned structured knowledge in RL training [11]. Based on the form of causal knowledge, we divide the related works into two categories, i.e., *implicit* and *explicit* causation. With *implicit* causal representation, researchers ignore the detailed causal structure. For instance, [12] extracts invariant features as one node that influences the reward function, while the other node consists of task-irrelevant features [13, 14, 15, 16]. This neat structure has good scalability but requires access to multiple environments that share the same invariant feature [12, 16, 17]. One can also turn to the *explicit* side by estimating detailed causal structures [18, 19, 20, 21],

Figure 1: The agent picks and places objects to achieve given goals. **Left:** The causal graph of this task. **Right:** Training setting: the hexagon is always pink, and the box is always blue. Three testing settings: (1) In distribution: the same as the training setting. (2) Spuriousness: swap the color and shape to break the spurious correlation. (3) Composition: increase the height of the goal.

which uses directed graphical models to capture the causality in the environment. A pre-request for this estimation is the object-level or event-level abstraction of the observation, which is available in most tasks and also becoming a frequently studied problem [22, 23, 24]. However, existing *explicit* causal reasoning RL models either rely on heuristic design without theoretical guarantees [18] or require the true causal graph [25].

In this paper, we propose *GeneRAlizing by DiscovERing (GRADER)*, a causal reasoning method that augments the RL algorithm with data efficiency, interpretability, and generalizability. We mainly focus on Goal-Conditioned RL (GCRL), where different goal distributions during training and testing reflect the generalization. We formulate the GCRL into a probabilistic inference problem [26] with a learnable causal graph as the latent variable. This novel formulation naturally explains the learning objective with three components, i.e., transition model learning, planning, and causal graph discovery, leading to an optimization framework that alternates between the causal structure discovery policy learning to gain generalizability. Under some mild conditions, we prove the unique identifiability of the causal graph and the theoretical performance guarantee of the proposed framework.

To demonstrate the effectiveness of the proposed method, we conduct comprehensive experiments in environments that require strong reasoning capability. Specifically, we design two types of generalization settings, i.e., spuriousness and composition, and provide an example to illustrate these settings in Figure 1. The evaluation results confirm our method's advantages in several aspects. First, the proposed data-efficient discovery method provides an explainable causal graph yet requires much less data than score-based methods, increasing data efficiency and interpretability during task solving. Second, simultaneously discovering the causal graph during policy learning dramatically increases the success rate of solving tasks. In summary, the contribution of this paper is threefold:

- We use the causal graph as a latent variable to reformulate the GCRL problem and then derive an iterative training framework from solving this problem.

- We prove that our method uniquely identifies true causal graphs, and the performance of the iterative optimization is guaranteed with a lower bound given converged transition dynamics.

- We design nine tasks in three environments that require strong reasoning capability and show the effectiveness of the proposed method against five baselines on these tasks.

2 Problem Formulation and Preliminary

We start by discussing the setting of RL we use in this paper and the assumptions required in causal reasoning. Then we briefly introduce the necessary concepts related to causality and causal discovery.

2.1 Factorized Goal-conditioned RL

We assume the environment follows the Goal-conditioned Markov Decision Process (MDP) setting with full observation. This setting is represented by a tuple $\mathcal{M} = (\mathcal{S}, \mathcal{A}, \mathcal{P}, \mathcal{R}, G)$, where $\mathcal{S}$ is the state space, $\mathcal{A}$ is the action space, $\mathcal{P}$ is the probabilistic transition model, $G \subset \mathcal{S}$ is the goal space which is a set of assignment of values to states, and $r(s, g) = \mathbb{1}(s = g) \in \mathcal{R}$ is the sparse deterministic reward function that returns 1 only if the state s match the goal g. In this paper, we focus on the goal-conditioned generalization problem, where the goal for training and testing stages will be sampled from different distributions $p_{\text{train}}(g)$ and $p_{\text{test}}(g)$. We refer to a goal $g \in G$ as a task and use these two terms interchangeably. To accomplish the causal discovery methods, we make a further assumption similar to [20, 27] for the state and action space:

Assumption 1 (Space Factorization). *The state space $\mathcal{S} = \{\mathcal{S}_1 \times \cdots \times \mathcal{S}_M\}$ and action space $\mathcal{A} = \{\mathcal{A}_1 \times \cdots \times \mathcal{A}_N\}$ can be factorized to disjoint components $\{\mathcal{S}_i\}_{i=1}^{M}$ and $\{\mathcal{A}_i\}_{i=1}^{N}$.*

The components representing one event or object's property usually have explicit semantic meanings for better interpretability. This assumption can be satisfied by state and action abstraction, which has been widely investigated in [22, 23, 24]. Such factorization also helps deal with the high-dimensional states since it might be impossible to treat each dimension as one random variable [18].

2.2 Causal Reasoning with Graphical Models

Reasoning with causality relies on specific causal structures, which is commonly represented as directed acyclic graphs (DAGs) [28] over variables. Consider random variables $\boldsymbol{X} = (X_1, \ldots, X_d)$ with index set $\boldsymbol{V} := \{1, \ldots, d\}$. A graph $\mathcal{G} = (\boldsymbol{V}, \mathcal{E})$ consists of nodes $\boldsymbol{V}$ and edges $\mathcal{E} \subseteq \boldsymbol{V}^2$ with (v, v) for any $v \in \boldsymbol{V}$. A node i is called a parent of j if $e_{ij} \in \mathcal{E}$ and $e_{ji} \notin \mathcal{E}$. The set of parents of j is denoted by $\mathbf{PA}_j^{\mathcal{G}}$. We formally discuss the graph representation of causality with the definitions:

Definition 1 (Structural Causal Models [28]). *A structural causal model (SCM) $\mathfrak{C} := (\boldsymbol{S}, \boldsymbol{U})$ consists of a collection $\boldsymbol{S}$ of d functions,*

$$X_j := f_j(\mathbf{PA}_j^{\mathcal{G}}, U_j), \quad j \in [d], \tag{1}$$

where $\mathbf{PA}_j \subset \{X_1, \ldots, X_d\} \backslash \{X_j\}$ are called parents of X_j; and a joint distribution $\boldsymbol{U} = \{U_1, \ldots, U_d\}$ over the noise variables, which are required to be jointly independent.

Definition 2 (Causal Graph [28], CG). *The causal graph $\mathcal{G}$ of a SCM is obtained by creating one node for each X_j and drawing directed edges from each parent in $\mathbf{PA}_j^{\mathcal{G}}$ to X_j.*

We note that CG describes the structure of the causality, and SCM further considers the specific causation from the parents of X_j to X_j as well as exogenous noises U_j. To uncover the causal structure from data distribution, we assume that the CG satisfies the *Markov Property* and *Faithfulness* ([28]), which makes the independences consistent between the joint distribution $P(X_1, \ldots, X_n)$ and the graph $\mathcal{G}$. We also follow the *Causal Sufficiency* assumption [29] that supposes we have measured all the common causes of the measured variables.

Existing work [30, 20] believes that two objects have causation only if they are close enough while there is no edge between them if the distance is large. Instead of using such a local view of the causality, we assume the causal graph is consistent across all time steps. Still the causal influence could be local, and the specific influence indicated by edges is estimated by the function $f_j(\mathbf{PA}_j, U_j)$.

3 Generalizing by Discovering (GRADER)

With proper definitions and assumptions, we now look at the proposed method. We first derive the framework of GRADER by formulating the GCRL problem as a latent variable model, which provides a variational lower bound to optimize. Then, we divide this objective function into three parts and iteratively update them with a specific performance guarantee.

Figure 2: The paradigm of GRADER.

3.1 GCRL as Latent Variable Models

The general objective of RL is to maximize expected reward function w.r.t. a learnable policy model π. Particularly, in the goal-conditioned setting, such objective is represented as $\max_\pi \mathbb{E}_{\tau \sim \pi, g \sim p(g)}[\sum_{t=0}^{T} r(s^t, g)]$, where $p(g)$ is the distribution of goals and $\tau := \{s^0, a^0, \ldots, s^T\}$ the action-state trajectory with maximal time step T. The trajectory ends only if the goal is achieved $g = s^T$ or the maximal step is reached.

Inspired by viewing the RL problem as probabilistic inference [31, 26], we replace the question from *"How to find actions to achieve the goal?"* to *"what are the actions if we achieve the goal?"*, leading to a likelihood maximization problem for $p(\tau|s^*)$ with $s^* := \mathbb{1}(g = s^T)$. Different from previous

work [32] that recasts actions as latent variables and inferring actions that result in "observed" high reward, we decompose $p(\tau|s^*)$ with $\mathcal{G}$ as the latent variable to get the evidence lower bound (ELBO)

$$\log p(\tau|s^*) = \log \int p(\tau|\mathcal{G}, s^*)p(\mathcal{G}|s^*)d\mathcal{G} \geq \mathbb{E}_{q(\mathcal{G}|\tau)}[\log p(\tau|\mathcal{G}, s^*)] - \mathbb{D}_{\mathrm{KL}}[q(\mathcal{G}|\tau)||p(\mathcal{G})], \quad (2)$$

where the prior $p(\mathcal{G})$ and variational posterior $q(\mathcal{G}|\tau)$ represent distributions over a graph structure, i.e, the probability of the existences of edges in the graph. $\mathbb{D}_{\mathrm{KL}}$ represents the Kullback–Leibler (KL) divergence between two graphs, which will be explained in Section 3.3. Recall that the space of goal G is a subset of the state, here we extend the meaning of g and assume all trajectories achieve the goal in the final state [33], i.e., $g = s^T$. Such extension makes it possible to further decompose the first term of (2) as (refer to Appendix A.1.2)

$$\log p(\tau|\mathcal{G}, s^*) = \log p(s^0) + \sum_{t=0}^{T-1} \log p(s^{t+1}|s^t, a^t, \mathcal{G}) + \sum_{t=0}^{T-1} \log \pi(a^t|s^t, s^*, \mathcal{G}) + \log p(g). \quad (3)$$

Here, we use the fact that $\mathcal{G}$ is effective in both the transition model $p(s^{t+1}|a^t, s^t, \mathcal{G})$ and the policy model $\log p(a^t|s^t, s^*, \mathcal{G})$, g only influences the policy model, and the initial state s_0 depends neither on $\mathcal{G}$ nor g. We assume that both initial state $\log p(s_0)$ and goal $\log p(g)$ follow the uniform distribution; thus, the first and last terms of (3) are constants. The policy term π, selecting action a^t according to both current state s^t and goal g, is implemented with planning method and is further discussed in Section 3.2. Finally, we maximize the likelihood $p(\tau|s^*)$ with the following reformulated ELBO as the objective

$$\mathcal{J}(\theta, \phi) = \mathbb{E}_{q_\phi(\mathcal{G}|\tau)} \sum_{t=0}^{T-1} \left[\log p_\theta(s^{t+1}|s^t, a^t, \mathcal{G}) + \log \pi_\theta(a^t|s^t, s^*, \mathcal{G}) \right] - \mathbb{D}_{\mathrm{KL}}[q_\phi(\mathcal{G}|\tau)||p(\mathcal{G})] \quad (4)$$

where θ is the shared parameter of transition model $p_\theta(s^{t+1}|a^t, s^t, \mathcal{G})$ and policy $\pi_\theta(a^t|s^t, s^*, \mathcal{G})$, and ϕ is the parameter of causal graph $q_\phi(\mathcal{G}|\tau)$. To efficiently solve this optimization problem, we iteratively updates parameter ϕ (causal discovery, Section 3.3) and parameter θ (model and policy learning, Section 3.2), as shown in Figure 2. Intuitively, these processes can be respectively viewed as the discovery of CG and update of SCM, which share tight connections discussed in Section 2.2.

3.2 Model and Policy Learning

Let us start with a simple case where we already obtain a $\mathcal{G}$ and use it to guide the learning of parameter θ via $\max_\theta \mathcal{J}(\theta, \phi)$. Since the KL divergence of $\mathcal{J}(\theta, \phi)$ does not involve θ, we only need to deal with the first expectation term, the likelihood of the transition model and policy. For the transition $p_\theta(s^{t+1}|a^t, s^t, \mathcal{G})$, we connect it with causal structure by further defining a particular type of CG and denote it as $\mathcal{G}$ in the rest of this paper:

Definition 3 (Transition Causal Graph). *We define a bipartite graph $\mathcal{G}$, whose vertices are divided into two disjoint sets $\mathcal{U} = \{\mathcal{A}^t, \mathcal{S}^t\}$ and $\mathcal{V} = \{\mathcal{S}^{t+1}\}$. $\mathcal{A}^t$ represents action nodes at step t, $\mathcal{S}$ state nodes at step t, and $\mathcal{S}^{t+1}$ the state nodes at step $t + 1$. All edges start from set $\mathcal{U}$ and end in set $\mathcal{V}$.*

Model learning. This definition builds the causal graph between two consecutive time steps, which indicates that the values of states in step $t + 1$ depend on values in step t. It also implies that the interventions [28] on nodes in $\mathcal{U}$ are directly obtained since they have no parents. We denote the marginal distribution of $\mathcal{S}$ as $p_{\mathcal{I}_\pi^s}$, which is collected by RL policy π. Combined with the Definition 1 of SCM, we find that $p_\theta(s^{t+1}|a^t, s^t, \mathcal{G})$ essentially approximates a collection of functions f_j following the structure $\mathcal{G}$, which take as input the values of parents of the state node s_j and outputs the value s_j. Thus, we propose to model the transition corresponding to $\mathcal{G}$ with a collection of neural networks $f_\theta(\mathcal{G}) := \{f_{\theta_j}\}_{j=1}^M$ to obtain

$$s_j^{t+1} = f_{\theta_j}([\mathbf{PA}_j^\mathcal{G}]^t, U_j), \quad (5)$$

where $[\mathbf{PA}_j^\mathcal{G}]^t$ represents the values of all parents of node s_j^t at time step t and U_j follows Gaussian noise $U_j \sim \mathcal{N}(0, \mathbf{I})$. In practice, we use Gated Recurrent Unit [34] as f_j because it supports varying numbers of input nodes. We take s_j^t as the initial hidden embedding and the rest parents $[\mathbf{PA}_j^\mathcal{G} \backslash s_j]^t$ as the input sequence to f_j. The entire model is optimized by stochastic gradient descent with the log-likelihood $\log p_\theta(s^{t+1}|a^t, s^t, \mathcal{G})$ as objective.

167 **Policy learning with planning.** Then we turn to the policy term $\pi_\theta(a^t|s^t, s^*, \mathcal{G})$ in $\mathcal{J}(\theta, \phi)$. We
168 optimize it with planning methods that leverage the estimated transition model. Specifically, the policy
169 aims to optimize an action-state value function $Q(s^t, a^t) = \mathbb{E}\left[\sum_{t'=0}^{H} \gamma^{t'} r(s^{t'+t}, a^{t'+t})|s^t, a^t\right]$,
170 which can be obtained by unrolling the transition model with a horizon of H steps and discount
171 factor γ. In practice, we use model predictive control (MPC) [35] with random shooting [36], which
172 selects the first action in the fixed-horizon trajectory that has the highest action-state value $Q(s^t, a^t)$,
173 i.e. $\hat{\pi}(s^t) = \arg\max_{a^t \in \mathcal{A}} Q_\theta^{\mathcal{G}}(s^t, a^t)$. The formulation we derived so far is highly correlated to the
174 model-based RL framework [37]. However, the main difference is that we obtain it with variational
175 inference by regarding the causal graph as a latent variable.

3.3 Data-Efficient Causal Discovery

177 In this step, we relax the assumption of knowing $\mathcal{G}$
178 and aim to estimate the posterior distribution $q(\mathcal{G}|\tau)$
179 to increase the ELBO (4) w.r.t. parameter ϕ. In most
180 score-based methods [38], likelihood is used to evaluate
181 the correctness of the causal graph, i.e., a better causal
182 graph leads to a higher likelihood. Since the first term
183 of (4) represents the likelihood of the transition model,
184 we convert the problem of $\max_\phi \mathcal{J}(\theta, \phi)$ to the causal
185 discovery that finds the true causal graph based on
186 collected data samples. As for the second term of (4),
187 the following proposition shows that the KL divergence
188 between $q_\phi(\mathcal{G}|\tau)$ and $p(\mathcal{G})$ can be approximated by a
189 sparsity regularization (proof in Appendix A.4.2).

190 **Proposition 1** (KL Divergence as Sparsity Regular-
191 ization)**.** *With entry-wise independent Bernoulli prior*
192 *$p(\mathcal{G})$ and point mass variational distribution $q(\mathcal{G}|\tau)$ of*
193 *DAGs, $\mathbb{D}_{KL}[q_\phi\|p]$ is equivalent to an ℓ_1 sparsity regu-*
194 *larization for the discovered causal graph.*

Algorithm 1: GRADER Training

Input: Trajectory buffer $\mathcal{B}_\tau$, Causal
   graph $\mathcal{G}$, Transition model f_θ
while θ *not converged* **do**
  // Policy from planning
  Sample a goal $g \sim p_{\text{train}}(g)$
  while $t < T$ **do**
    $a^t \leftarrow$ Planner(f_θ, s^t, g)
    $s^{t+1}, r^t \leftarrow$ Env(a^t, g)
    $\mathcal{B}_\tau \leftarrow \mathcal{B}_\tau \cup \{a^t, s^t, s^{t+1}\}$
  // Estimate causal graph
  for $i \leq M + N$ **do**
    for $j \leq M$ **do**
      Infer edge $e_{ij} \leftarrow q_\phi(\cdot|\mathcal{B}, \eta)$
  // Learn transition model
  Update $f_\theta(\mathcal{G})$ via (5) with $\mathcal{B}$

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195 We restrict the posterior $q_\phi(\mathcal{G}|\tau)$ to point mass distribution and use a threshold η to control the sparsity.
196 We perform the discovery process from the classification perspective by proposing binary classifiers
197 $q_\phi(e_{ij}|\tau, \eta)$ to determine the existence of an edge e_{ij}. This classifier $q_\phi(e_{ij}|\tau, \eta)$ is implemented by
198 statistic *Independent Test* [39], and η is the threshold for the p-value of the hypothesis. A larger η
199 corresponds to harder sparsity constraints, thus leading to a sparse $\mathcal{G}$ since two nodes are more likely
200 to be considered independent. According to the definition 3, we only need to conduct classification to
201 edges connecting nodes between $\mathcal{U}$ and $\mathcal{V}$. If two nodes are dependent, we add one edge directed from
202 the node in $\mathcal{U}$ to the node in $\mathcal{V}$. This definition also ensures that we always have $q(\mathcal{G}|\tau) \in \mathcal{Q}_{DAG}$,
203 where $\mathcal{Q}_{DAG}$ is the class of DAG. With this procedure, we identify a unique CG $\mathcal{G}^*$ under optimality:

204 **Proposition 2** (Identifiability)**.** *Given an oracle independent test, with an optimal interventional data*
205 *distribution $p^*_{\mathcal{T}^s_\pi}$, causal discovery obtains ϕ^* that correctly tells the independence between any two*
206 *nodes, then the causal graph is uniquely identifiable, with $e^*_{ij} = q_{\phi^*}(e_{ij}|\tau), \forall i \in [M+N], j \in [M]$.*

207 In practice, we use χ^2-test for discrete variables and the Fast Conditional Independent Test [39]
208 for continuous variables. The sample size needed for obtaining the oracle test has been widely
209 investigated [40]. However, testing with finite data is not a trivial problem, as stated in [41], especially
210 when the data is sampled from a Goal-conditioned MDP. Usually, the random policy is not enough
211 to satisfy the oracle assumption because some nodes cannot be fully explored when the task is
212 complicated and has a long horizon. To make this assumption empirically possible, it is necessary to
213 simultaneously optimize $\pi_\theta(a^t|s^t, s^*, \mathcal{G})$ to access more samples close to finishing the task, which is
214 further analyzed in Section 3.4. We also empirically support this argument in Section 4.2 and provide
215 a detailed theoretical proof in Appendix A.3 and A.4.

3.4 Analysis of Performance Guarantee

217 The entire pipeline of GRADER is summarized in Algorithm 1. To analyze the performance of the
218 optimization of (4), we first list important lemmas that connect previous steps and then show that the
219 iteration of these steps in *GRADER* leads to a virtuous cycle.

By the following lemmas, we show the following performance guarantees step by step. Lemma 1 shows model learning is monotonically better at convergence given a better causal graph from causal discovery. Then the learned transition model helps to improve the lower bound of value function during planning according to Lemma 2. Lemma 3 reveals the connection between policy learning and interventional data distribution, which in turn improves the quality of our causal discovery, as is shown in Lemma 4 and Proposition 2.

Lemma 1 (Monotonicity of Transition Likelihood). *Assume* $\mathcal{G}^* = (V, E^*)$ *be the true CG, for two CG* $\mathcal{G}_1 = (V, E_1)$ *and* $\mathcal{G}_2 = (V, E_2)$, *if* $SHD(\mathcal{G}_1, \mathcal{G}^*) < SHD(\mathcal{G}_2, \mathcal{G}^*)$, $\exists \, e, \, s.t. \, E_1 \cup \{e\} = E_2$, *when transition model* θ *converges, the following inequality holds for the transition model in (4):*

$$\log p_\theta(s^{t+1}|a^t, s^t, \mathcal{G}^*) \geq \log p_\theta(s^{t+1}|a^t, s^t, \mathcal{G}_1) \geq \log p_\theta(s^{t+1}|a^t, s^t, \mathcal{G}_2) \tag{6}$$

where SHD is the Structural Hamming Distance defined in Appendix A.2.

Lemma 2 (Bounded Value Function in Policy Learning). *Given a planning horizon* $H \to \infty$, *if we already have an approximate transition model* $\mathbb{D}_{TV}(\hat{p}(s'|s, a), p(s'|s, a)) \leq \epsilon_m$, *the approximate policy* $\hat{\pi}$ *achieves a near-optimal value function (refer to Appendix A.3 for detailed analysis):*

$$\|V^{\pi^*}(s) - V^{\hat{\pi}}(s)\|_\infty \leq \frac{\gamma}{(1-\gamma)^2}\epsilon_m \tag{7}$$

Lemma 3 (Policy Learning Improves Interventional Data Distribution). *With a step reward* $r(s, a) = \mathbb{1}(s = g)$, *we show that the value function determines an upper bound for TV divergence between the interventional distribution with its optimal (proof details in Appendix A.3):*

$$\mathbb{D}_{TV}(p_{\mathcal{I}_\pi^s}, p_g) \leq 1 - (1 - \gamma)V^\pi(s). \tag{8}$$

where $p_{\mathcal{I}_\pi^s}$ *is the marginal state distribution in interventional data, and* p_g *is the goal distribution. A better policy with larger* $V^\pi(s)$ *enforces the distribution of interventional data towards the goal.*

Lemma 4 (Interventional Data Benefits Causal Discovery). *For* $\epsilon_g = \min_{p_g>0} p_g$, $\mathbb{D}_{TV}(p_{\mathcal{I}_\pi^s}, p_g) < \epsilon_g$, *the error of our causal discovery is upper bounded with* $\mathbb{E}_{\hat{\mathcal{G}}}[SHD(\hat{\mathcal{G}}, \mathcal{G}^*)] \leq |\mathcal{S}| - 1$.

After the close-loop analysis of our model, we are now able to analyze the overall performance of the proposed framework. Under the construction of $p_\theta(s^{t+1}|a^t, s^t, \mathcal{G})$ with NN-parameterized functions, the following theorem shows that the learning process will guarantee to perform a close estimation of true ELBO under the iterative optimization among model learning, planning, and causal discovery.

Theorem 1. *With T-step approximate transition dynamics* $\mathbb{D}_{TV}\Big(\hat{p}(s'|s, a), p(s'|s, a)\Big) \leq \epsilon_m$, *if the goal distribution satisfies* $\epsilon_g > \frac{\gamma}{1-\gamma}\epsilon_m$, *and the distribution prior CG is entry-wise independent Bernoulli($\epsilon_\mathcal{G}$), GRADER guarantees to achieve an approximate ELBO* $\hat{\mathcal{J}}$ *with the true ELBO* $\mathcal{J}^*$:

$$\|\mathcal{J}^*(\theta, \phi) - \hat{\mathcal{J}}(\hat{\theta}, \hat{\phi})\|_\infty \leq \left[1 + \frac{\gamma}{(1-\gamma)^2}\right]\epsilon_m T + \log\left(\frac{1 - \epsilon_\mathcal{G}}{\epsilon_\mathcal{G}}\right)(|\mathcal{S}| - 1), \tag{9}$$

An intuitive understanding of the performance guarantee is that a better transition model indicates a better approximation of objective $\mathcal{J}$. Yet it's harder for *GRADER* to approximate a longer trajectory with large state space. The proof of this theorem and corresponding empirical results are in Appendix A.5.

4 Experiments

In this section, we first discuss the setting of our designed environments as well as the baselines used in the experiments. Then, we provide the numerical results and detailed discussions to answer the following important research questions: **Q1.** Compared to baselines, how does GRADER gain performance improvement under both in-distribution and generalization settings? **Q2.** Compared to an offline random policy, how does a well-trained policy improve the results of causal discovery? **Q3.** Compared to score-based causal discovery, does the proposed data-efficient causal discovery pipeline guarantee to identifying true causal graph as stated in Section 3.3? **Q4.** Compared to false causal graphs, how does the correctness of discovered causal graph influence the task-solving performance of RL agents?

(1) Stack   (2) Unlock   (3) Crash

Figure 3: Three environments used in this paper.

Table 1: Success rate (%) for nine settings in three environments. **Bold** font means the best.

Method	Stack-I	Stack-S	Stack-C	Unlock-I	Unlock-S	Unlock-C	Crash-I	Crash-S	Crash-C
SAC	34.7±16.1	22.1±14.0	31.7±5.1	0.1±0.5	0.0±0.2	0.4±1.7	22.5±17.6	18.6±8.7	6.7±3.8
ICIN	71.8±6.9	71.0±7.4	58.6±8.3	31.7±9.6	32.7±8.6	31.5±8.5	27.9±6.1	15.8±17.2	7.8±8.8
PETS	**97.2**±**6.9**	77.7±13.5	73.7±10.3	59.5±7.2	20.6±5.9	28.3±10.0	52.3±11.5	44.6±12.5	37.1±5.1
TICSA	85.9±8.4	88.8±10.1	76.2±8.3	58.5±12.3	33.6±14.3	29.8±8.3	68.9±5.9	56.8±8.6	15.0±8.2
ICIL	93.7±5.9	81.2±14.4	62.8±13.0	**67.1**±**11.6**	15.9±4.7	53.6±15.3	55.3±20.9	21.7±17.7	14.3±7.3
GNN	45.7±9.1	39.0±10.4	41.7±8.6	3.4±2.3	3.4±2.4	4.5±3.0	4.2±4.0	5.1±5.1	3.8±2.8
GRADER	95.6±5.4	**97.6**±**6.0**	**93.7**±**8.4**	64.2±9.1	**61.4**±**4.4**	**82.1**±**9.2**	**91.5**±**4.4**	**84.3**±**10.0**	**84.7**±**7.3**
Score	92.7±7.4	90.5±7.5	73.9±8.5	44.9±28.1	23.1±7.6	36.2±30.1	42.3±17.5	53.4±18.7	8.4±6.1
Full	92.9±6.3	86.0±9.5	75.7±10.3	63.8±9.2	18.3±7.4	53.7±14.3	69.8±14.0	52.6±12.8	42.0±17.2
Offline	96.8±5.8	95.4±6.1	81.4±7.8	13.8±8.1	13.9±7.5	11.7±6.9	13.1±16.2	30.2±16.5	14.9±12.4

4.1 Environments and Baselines

Since most commonly used RL benchmarks do not explicitly require causal reasoning for generalization, we design three environments shown in Figure 3. These environments use the true state as observation to disentangle the reasoning task from visual understanding. For each environment, we design three settings – in-distribution (*I*), spuriousness (*S*), and composition (*C*) – corresponding to different goal distributions for generalization. We use $p_{\text{train}}(g)$ and $p_{\text{test}}(g)$ to represent the goal distribution during training and testing, respectively. *I* uses the same $p_{\text{train}}(g)$ and $p_{\text{test}}(g)$, *S* introduces spurious correlations in $p_{\text{train}}(g)$ but remove them in $p_{\text{test}}(g)$, and *C* contains more similar sub-goals in $p_{\text{test}}(g)$ than in $p_{\text{train}}(g)$. The details of these settings in three environments are briefly summarized in the following (details in in Appendix C.2.1):

- **Stack**: We design this manipulation task inspired by the CausalWorld [42], where the agent must stack objects to match specific shapes and colors. In *Stack-S*, we let the same shape have the same color in $p_{\text{train}}(g)$ but randomly sample the color and shape in $p_{\text{test}}(g)$. In *Stack-C*, the maximum number of object is two in $p_{\text{train}}(g)$ but five in $p_{\text{test}}(g)$.

- **Unlock**: We design this task for the agent to collect a key to open doors built upon the Minigrid [43]. In *Stack-S*, the door and the key are always in the same row in $p_{\text{train}}(g)$ but uniformly sample in $p_{\text{test}}(g)$. In *Unlock-C*, there are one door in $p_{\text{train}}(g)$ but two doors in $p_{\text{test}}(g)$.

- **Crash**: The occurrence of accidents usually relies on causality, e.g., an autonomous vehicle (AV) collides with a jaywalker because its view is blocked by another car [44]. We design such a crash scenario based on highway-env [45], where the goals are to create crashes between a pedestrian and different AVs. In *Stack-S*, the initial distance between ego car and pedestrian is a constant in $p_{\text{train}}(g)$ but irrelevant in $p_{\text{test}}(g)$. In *Stack-C*, there is one pedestrian in $p_{\text{train}}(g)$ but two in $p_{\text{test}}(g)$.

We use the following methods as our baselines to fairly demonstrate the advantages of GRADER. **SAC:** [46] Soft Actor-Critic is a well-known model-free RL method; we augment it with HER [47] in *Stack* because the goal in this task does not have pre-conditions. **ICIN:** [25] It uses DAgger [48] to learn goal-conditioned policy with the causal graph estimated from the expert policy. We assume it can access the true causal graph for supervised learning. **PETS:** [49] We consider the ensemble transition model with random shoot planning as one baseline, which achieves generalization with the uncertainty-aware design. **TICSA:** [18] This is a causal-augmented MBRL method that simultaneously optimizes a soft adjacent matrix (representing the causality) and a transition model. **ICIL:** [16] This method proposes an invariant feature learning structure that captures the implicit causality of multiple tasks. We only use it for transition model learning since the original method is designed for imitation learning. **GNN:** [50] Since graph neural network (GNN) is good at learning structural information, we implement a GNN-based baseline using Relational Graph Convolutional Network.

4.2 Results Discussion

Overall Performance (Q1) We compare the testing reward of all methods under nine tasks and summarize the results in Table 1 to demonstrate the overall performance. Generally, our method outperforms other baselines in all tasks except *Stack-I*, which is simple for all methods. We note that the gap between our method and baselines under *S* and *C* settings is more significant than the *I* setting, showing that our method still works well in the non-trivial generalization task. As a model-free

Figure 4: **Left:** Test reward of the *Crash* environment calculated by 30 trails. **Right:** The accuracy of causal graph discovery with samples from GRADER, Score, and Offline (random policy).

method, SAC fails in all three tasks of *Unlock* and *Crash* environments since they have sparse rewards. Without learning the causal structure of the environment, PETS even cannot fully solve *Unlock-S*, *Unlock-C*, and all *Crash* tasks. Both TICSA and ICIL learn explicit or implicit causality of the task, so they are relatively better than SAC and PETS. However, they are still worse than GRADER in two generalization settings because of the unstable and inefficient causal reasoning mechanism. We also find that even if the ICIN is given the true causal graph, the policy learning part cannot efficiently leverage the causality, leading to worse performance in generalization settings.

To further analyze the tendency of learning, we plot the curves of all methods under *Crash* in Figure 4. Our method quickly learns to solve the tasks at the beginning of the training, demonstrating data efficiency. GRADER also outperforms other methods with large gaps in the later training phase. The results of the other two environments can be found in Appendix B.1.

Importance of Policy Learning (Q2) As we mentioned in Section 3.3, we empirically compare GRADER and **Offline** [51], which uses data from offline random policy, and plot results in the right part of Figure 4. We use SHD [52] to compute the distance between the estimated causal graph and the true causal graph, which is obtained by checking the best causal graph with human analysis. The true causal graph for each environment can be found in Appendix B.2. When we only use samples collected offline by random policy, we cannot obtain variables' values that require long-horizon reasoning, e.g., the door can be opened if the agent is close to the door and have the key, leading to unreliable discovery. As a consequence, the causal graph obtained by Offline indeed harms the performance, as shown in Figure 4. Instead, GRADER gradually explores more regions and quickly obtains the true causal graph when we iteratively discover the causal graph and update the policy.

Advantage of Data-efficient Causal Discovery (Q3) To show the advantage of proposed constraint-based methods, we design a model named **Score** that optimizes a soft adjacent matrix similar to the score-based method in the discovery step. Score-based discovery methods, also used in TICSA, are recently combined with NN for differentiable training. According to the discovery accuracy shown in the right part of Figure 4, we find score-based discovery is inefficient. Based on the performance of the Score model summarized in Table 1, we also conclude that it is not as good as our constraint-based method and has a large variance due to the unstable learning of causal graph.

Figure 5: Influence of different causal graphs on *Unlock-S*.

Influence of Causal Graph (Q4) To illustrate the importance of the causal graph, we implement another variant of GRADER named **Full**, which uses a fixed full graph that connects all nodes between the sets $\mathcal{U}$ and $\mathcal{V}$. According to the performance shown in Table 1 and Figure 4, we find that the full graph achieves worse results than GRADER because of the redundant and spurious correlation. Intuitively, unrelated information causes additional noises to the learning procedure, and the spurious correlation creates a shortcut that makes the model extract wrong features, leading to worse results in the spuriousness generalization as shown in Table 1.

We then investigate how the correctness of the causal graph influences the performance. We use fixed graphs interpolating from the best causal graph to the full graph to train a GRADER model on

Unlock-S and summarize the results in Figure 5 The more correct the graph is, the higher reward the agent obtains, which supports our statements in Section 3.4 that the causal graph is important for the reasoning tasks – a better causal graph helps the model have better task-solving performance.

5   Related Works

RL Generalization From agent's view, algorithms focus on actively obtaining the structure of the task or environment. Some decompose the given task into sub-tasks [53, 54, 55, 5] and when they encounter an unseen task, they rearrange those sub-tasks to solve new tasks. Instead of dividing tasks, Symbolic RL learns a program-based policy consisting of domain language [6, 8] or grammar [7, 56]. The generated program is then executed [9] to interact with the environment, which has the potential to solve unseen tasks by searching the symbolic space. From the other view, we generate environments to augment agents' experience for better generalization. One straightforward way is data augmentation of image-based observation [57, 58, 59, 60, 61, 62]. When extended to other factors of the environment, *Domain Randomization* [63] and its variants [64, 65] are proposed.

Goal-Conditioned RL (GCRL) From the optimization perspective, Universal Value Function [66], reward shaping [67], and latent dynamic model [68] are widely used tools to solve GCRL problem. Sub-goal generation [69] is another intuitive idea to tackle the long horizon with sparse reward, where the core thing is to make sure the generated sub-goals are solvable. Finally, *HER* [33], belonging to the relabelling category, is a ground-breaking yet straightforward method that treats visited states as "fake" goals when the goal and state share the same space. Later on, improved versions of HER [47, 70, 71] were widely studied. One limitation is that we cannot directly use a visited state as a goal if the goal has pre-conditions. Similar to our setting, [72] and [73] convert the GCRL problem to variational inference by regarding control as inference [26]. [72] propose a EM framework under the HER setting and [73] treats the last state as goal and estimate a shaped reward during training.

RL with Causal Reasoning Causality is now frequently discussed in the machine learning field to complement the interpretability of neural networks [28]. RL algorithms also incorporate causality to improve the reasoning capability [74]. For instance, [25] and [19] explicitly estimate causal structures with the interventional data obtained from the environment. These structures are either combined with dynamic model learning in model-based RL [18] or used to adjust the buffer priority [20]. Implicitly, we can abstract the causal structure and formulate it using the *Block MDP* [12] setting or training multiple encoders to extract different kinds of representations [15]. Following the idea of invariant risk minimization [75], they assume task-relevant features are invariant and shared across all environments, which can be used as the only cause of the reward.

Causal Discovery Causal discovery [76] is a long-stand topic in economics and sociology, where the traditional methods can be generally categorized into constraint-based and score-based. Constraint-based methods [29] start from a complete graph and iteratively remove edges with conditional independent test [77, 78] as constraints. Score-based methods [38, 79] use metrics such as *Bayesian Information Criterion* [80] as scores and prefer edges that maximize the score given the dataset. Recently, researchers extend score-based methods with RL [81] or differentiable discovery [82, 83]. The former selects edges with a learned policy, and the latter learns a soft adjacent matrix with observational or interventional data. Active intervention methods are also explored [84] to increase the efficiency of data collection and decrease the cost of conducting intervention [85].

6   Conclusion

This paper proposes a latent variable model that injects a causal graph reasoning process into transition model learning and planning to solve GCRL problems under the generalization setting. We theoretically prove that our iterative optimization process can obtain the true causal graph. To evaluate the performance of the proposed method, we designed nine tasks in three environments. The comprehensive experiment results show that our method has better data efficiency and performance than baselines. Our method also provides interpretability by the explicitly discovered causal graph. The main limitation of this work is that the explicit estimation of causal structure does not scale well to the number of nodes. Developing efficient gradient-based discovery methods could be a promising direction. In addition, the factorized state and action space assumption may restrict the usage of this work to semantic representations, which need to be processed with abstraction methods. We further discuss the potential negative social impact and addition limitation in Appendix C.3.

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