## Multi-Agent Local Voting Protocol For Online DAG Scheduling

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# IC (O) MIP

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## Graph Scheduling / Distributed Heterogeneous Computing

- **Directed acyclic graph (DAG)**  $G = (J, E)$ ;  $J$  set of nodes,  $E$  set of edges:
- Node  $v \in J$  represent a computational task (job) with a given type of executor that it should be executed upon
- Edge  $(v_i, v_j) \in E$  denotes precedence constraint, i.e. node  $v_j$  cannot be executed until  $v_i$  and all its other predecessors are not complete
- Node without predecessors entry node, without successors exit node; there can be multiple of both types in a given DAG
- Cost of communication is  $b_{i,j}$  defined for all edges  $(v_i, v_j) \in E$  and should accounted for if  $v_i$  and  $v_j$  are executed by different executors

**Online sliding informational window**: at each given moment the **scheduler observes** only **ready-to-execute** nodes and their **immediate successors** within the **informational window with length 2**. This limits amount of information and makes problem online: **graph is not known in advance**, opposite to offline case

At each moment of time  $t$ , executors calculate a probability of assignment of a given task (node). Probability is calculated as a ratio of a given node's **aggregated metric** to sum of all other observable nodes **aggregated metrics.** It is constructed from following characteristics of a DAG and executors:

- **Incoming connections** ratio that represent how many unfinished parent nodes has each of a given node's child nodes  $W_t$
- **Available nodes –** ratio of tasks of a matching to the given executor's type that are ready-to-be-executed  $Q_i$
- **Relative performance difference –** difference between minimum execution time of a given node by all executors to a given executor  $Z_i$

Coefficients for sub-metrics are calculated using genetic algorithm with singlepoint crossover and random mutation, and selective fixed-size elitism. Assignment probabilities are synchronized using **Local Voting Protocol (LVP):** Each executor shares its task assignment probabilities with neighboring executors that iteratively update their task probabilities based on the state of their neighbors and **consensus** is reached when the difference in task assignment probabilities across all neighboring executors falls below a predefined threshold ( $\epsilon$  = 0.05)

**Distributed heterogeneous computing** consists of a set P where  $p_i \in P$  is heterogeneous executors with fully connected topology, for which defined:

- Computational cost matrix D, where  $D^{i,j} = w_{i,j}$  is the execution time required for executor  $p_i \in P$  to process node  $v_i \in J$
- Each executor has a defined type; to execute task  $v_j$  on  $p_i$  it is required that their types match

**Goal is to construct mapping of nodes from DAG to executors, such that** 

 $makespan \rightarrow min$ where makespan is max  $v \in V$ actual finish time of  $v$ ).

- **Executors:** 3, 6, 12, 24 (with  $n$  executors of each type 1, 3, 4, 8)
- **Number of nodes:** 3000, 6000, 12000, 24000 respectively to executors

**Use machine learning approaches** [3] to **predict computational tasks and dependencies between them** for accurate prediction of the next task's properties and its dependencies to the existing **task to extend online method's DAG knowledge and capabilities** Input Layer

[1] Multi-agent Reinforcement Learning-based Adaptive Heterogeneous DAG Scheduling, ACM Transactions on Intelligent Systems and Technology, Volume 14, Issue 503, 2023, pp 1–26 [2] Heterogeneous Computational Scheduling using Adaptive Neural Hyper-heuristic, Doklady Mathematics (accepted, will be printed next month) [3] DAGCN: hybrid model for efficiently handling joint node and link prediction i

Table 5 Similarity between Original and Predicted DAGs



#### Table 4 Prediction Quality



[Task Feature, liancency Matrix<sup>1</sup>



**MARL** [1] first level – make ordered list of tasks to be executed by each type of resource via Multi-Agent Proximal Policy Optimization with agent for each type

#### Online vs offline scheduling

#### MLVP



#### Example

DAGs generated DAGGEN [2] – industry standard for evaluating DAG heuristics



### Offline methods [1, 2]

- **State**: matrix of 13 DAG metrics [1] for each buffer of ready-to-execute node with pre-defined length (hyperparameter)
- **Action**: select metric by which order the list of ready-to-executed nodes buffer
- **Reward:** normalized difference between MILP and MARL makespan after each DAG scheduled (training on a batch of DAGs at once)
- **Second level –** allocate tasks from ordered list using Earliest Finish Time (EFT) **NN [2] First level –** make ordered list of tasks to be executed by each type:
- **Single agent :** Neural Network trained using Genetic Programming approach
- **Genetic programming allows to skip construction of differentiable makespan improvement function**
- **Features**: 8 DAG metrics [3] + type
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• **Architecture:** 3 fully connected layers + Genetic Programming algorithm **Second level –** allocate tasks from ordered list using Earliest Finish Time (EFT)



