# Efficient Execution of SPARQL Queries with OPTIONAL and UNION Expressions

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Abstract—The proliferation of RDF datasets has resulted in studies focusing on optimizing SPARQL query processing. Most existing work focuses on basic graph patterns (BGPs) and ignores other vital operators in SPARQL, such as UNION and OPTIONAL. SPARQL queries with these operators, which we abbreviate as SPARQL-UO, pose serious query planning challenges. In this paper, we propose techniques for optimizing SPARQL-UO queries using BGP execution as a building block, based on a novel <u>BGP-based Evaluation (BE)-Tree representa-</u> tion of query plans. On top of this, we propose a series of *cost-driven BE-tree transformations* to generate more efficient plans by reducing the search space and intermediate result sizes, and a *candidate pruning* technique that further enhances efficiency at query time. Experiments confirm that our method outperforms the state-of-the-art by orders of magnitude.

Index Terms—graph database, graph query, query optimization, OPTIONAL expressions, UNION expressions

## I. INTRODUCTION

The proliferation of knowledge graphs has generated many RDF (Resource Description Framework) data management problems. RDF is the de-facto data model for knowledge graphs, where each edge is a triple of (subject, predicate, object). SPARQL has been the focus of a significant body of research as the standard language for accessing RDF datasets. Most of the existing work focus on basic graph pattern (BGP) execution, the basic building block of SPARQL. On the other hand, how to execute and optimize queries containing operators on graph patterns, such as UNION and OPTIONAL, has received much less attention.

UNION and OPTIONAL expressions are essential in SPARQL grammar. Firstly, RDF datasets are semi-structured and do not enforce a schema. The UNION operator is crucial in this case since it merges diversely expressed information. For example, in DBpedia [1], an open-domain knowledge graph extracted from Wikipedia, persons' names are represented using the predicate foaf:name or rdfs:label. Thus, to fully retrieve all the names of a group of persons (e.g., Presidents of the United States), it is necessary to use the UNION operator (Fig. 1(a)). Secondly, RDF datasets are commonly incomplete. Specifically, an entity may lack some attributes or relationships that most other entities of the same type have. In this case, the OPTIONAL operator is useful, since it allows attaching some attributes or relations as optional information. For example, the OPTIONAL query in Fig. 1(b) fetches all the presidents of the United States, along with other references to them that are not on the same Wikipedia page. Since not every president has multiple references in the database, the triple with the predicate owl:sameAs is enclosed in an OPTIONAL expression.

UNION and OPTIONAL expressions are widely used in real-world SPARQL workloads and are part of the SPARQL 1.1 specification. Recent empirical studies [2] show that UNION and OPTIONAL expressions occur in 25.10% and 31.72% of the valid queries from real SPARQL query logs across a diverse range of endpoints, respectively. In this paper, we address the efficient execution of SPARQL queries with UNION and OPTIONAL expressions, which we abbreviate as SPARQL-UO queries.

**Our Solution.** Since BGP has been well studied, it is desirable to build SPARQL-UO query optimization on a well-performing BGP engine. Therefore, we first propose a BGP-based query evaluation scheme. Specifically, we propose a *B*GP-based *E*valuation (BE)-tree query plan representation for the evaluation plan of SPARQL-UO queries, which is a tree structure with BGPs as leaves and graph pattern operators as internal nodes. However, if the BE-tree is evaluated as it is, some BGPs may generate large intermediate results. Therefore, we propose a BE-tree transformation method to generate a more efficient query plan, which is backed by relational algebra (RA). The BE-tree serves as an easily constructible and comprehensible intermediate form between a SPARQL query and its RA-based plan.

We introduce two types of transformations, *merge* and *inject*, that target UNION and OPTIONAL operators, respectively. These transformations expose opportunities for reducing the cost during the evaluation of BGPs, UNION and OPTIONAL operators while preserving query semantics, which are one-to-one mapped from the BE-trees to the RA-based plans for execution. Since there are many different ways to transform an RA-based plan, we devise a cost model that accounts for the cost of evaluating both BGPs and these operators, and choose the transformation that reduces the most cost. Because of the vast space of possible transformations, we propose a greedy strategy to determine

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Fig. 1: An Example Query with a UNION and OPTIONAL Clause.

the transformations step-by-step. The transformed plan is then evaluated by the BGP-based scheme, enhanced by the query-time optimization, *candidate pruning*, which prunes the search space of BGP evaluation on the fly.

To summarize, we make the following contributions:

- We propose a novel BE-tree representation for the evaluation plan of a SPARQL-UO query and design two RAbacked BE-tree transformation primitives, *merge* and *inject*, to generate more efficient SPARQL-UO query plans.
- We propose a cost model for SPARQL-UO queries and a cost-driven BE-tree transformation algorithm.
- 3) We design a query-time optimization called *candidate pruning* that augments the BGP-based query evaluation scheme by pruning the search space.
- 4) We conduct experiments on large-scale real and synthetic RDF datasets, which shows that our method outperforms existing techniques by orders of magnitude.

The rest of the paper is organized as follows. A brief review of related work is given in Sec. II. Sec. III provides the necessary preliminary information. Sec. IV presents the BE-tree plan representation, the transformations, and the query evaluation scheme. The cost-driven planning algorithm is proposed in Sec. V, and the query-time optimization is presented in Sec. VI. We experimentally evaluate our method in Sec. VII and conclude the paper in Sec. VIII.

# II. RELATED WORK

Although SPARQL query optimization has been extensively studied, most of the focus has been on evaluating BGPs [3], including graph-based approaches and relational approaches. Graph-based approaches include works on effective index strategies (e.g., gStore [4]) and join order optimization (e.g., WCOJ [5]). In contrast, relational approaches rely on a relational DBMS and consider RDF graphs as threecolumn tables or other complex table organizations [6], [7]. Processing SPARQL queries is then mapped to its relational counterparts, as done in Apache Jena [8] and Virtuoso [9]. Relational BGP optimization approaches focus primarily on efficient data organization (e.g., property table [8], vertical partitioning [10] and single table exhaustive indexing [11]). However, these BGP optimization techniques cannot optimize UNION and OPTIONAL, since the semantics of these operators are fundamentally different from joins. As explained in later sections, our solution relies on BGP evaluation as a basic building block, and our proposed

optimization techniques operate on a higher level than BGP evaluation techniques.

Works on optimizing SPARQL queries with UNION and OPTIONAL operators are quite scarce. LBR [3] proposes a query rewriting technique to reduce intermediate results of left-outer joins, the join semantics represented by OPTIONAL. To remove inconsistent variable bindings, LBR uses the *nullification* and *best-match* techniques previously studied in SQL left-outer joins [12], and proposes a semijoin strategy to prune the candidate results. However, it follows an execution strategy of two-pass semi-join scans on the join graph, which introduces additional overhead during query execution. In this paper, we propose a more comprehensive approach that handles both UNION and OPTIONAL. Experiments also demonstrate that our techniques significantly outperform LBR on OPTIONAL queries. Works on ontology-based query reformulation [13], [14], which target SPARQL queries with UNION, bear a conceptual similarity to ours in that they consider a space of possible rewritten queries and select the one with the lowest cost for execution. However, the rewriting rules and cost model that they use are different from ours, and they do not consider OPTIONAL.

There are a number of theoretical works on optimizing queries with OPTIONAL [15], [16], which propose semantics-preserving rewriting rules for such queries. Our method, which operates on two particular rules, is parallel to these works and may be extended to support them in a cost-based manner in the future.

Note that our techniques to optimize UNION expressions can also be applied to conjunctive relational queries with unions due to their semantic similarity. In fact, all SPARQL-UO queries can be equivalently mapped to SQL, but the mapping of OPTIONAL expressions involves subselects in SQL, so our techniques cannot be applied without major adaptations [17], [18]. We are aware of a recent demonstration [19] that proposes a *join pushing* technique on conjunctive queries with unions, which pushes the join condition into the unioned sets if a cost model deems it more efficient. However, no description of the employed cost model is provided, which renders experimental comparison impossible. Some works on relational algebra have focused on outerjoin optimization [20], which can be applied to our setting. However, to our knowledge, there has not been similar works that treat rewritings in a systematic cost-based manner and outperform well-established baseline RDF systems experimentally.

**III.** PRELIMINARIES

# **RDF** Dataset.

Definition 1 (RDF dataset): Let pairwise disjoint infinite sets *I*, *B*, and *L* denote IRI, blank nodes and literals, respectively. An RDF dataset *D* is a collection of triples  $D = \{t_1, t_2, ..., t_{|D|}\}$ , where each triple is a three-tuple  $t = \langle subject, property, object \rangle \in (I \cup B) \times I \times (I \cup B \cup L)$ . Tab. I is an example RDF dataset with seven triples.

SPARQL Query-Syntax. In the following, we present the query dialect that we target in this paper, which is a proper subset of SPARQL 1.1. Assume there is an infinite set V representing the variables that appear in the query. All variables differ from IRIs and literals by leading with a question mark (?), so the set V is disjoint with I and L. This work focuses on SELECT queries, which retrieve results by matching the graph pattern in the query with the dataset. A SELECT query is of the form "SELECT  $v_1 v_2 \dots v_k$  WHERE {...}", in which the SELECT clause represents the query header, and the WHERE clause represents the query body (Fig. 2(a)). The SELECT clause determines the projection variables that need to appear in the query results, and the WHERE clause gives the group graph pattern that needs to be matched over the RDF dataset, which may consists of many other types of graph patterns, defined as follows.

Definition 2 (Triple Pattern): A triple  $t \in (V \cup I) \times (V \cup I) \times (V \cup I \cup L)$  is a triple pattern.

Basic graph patterns (BGPs) are sets of triple patterns. In this paper, we focus on connected BGPs, i.e., BGPs that represent connected query subgraphs. For brevity, all mentions of BGPs refer to connected BGPs henceforth. To give a formal definition of BGPs, we first introduce the notion of *coalescability*.

*Definition 3 (Coalescable triple patterns):* We say that the triple patterns  $t_1 = \langle s_1, p_1, o_1 \rangle$  and  $t_2 = \langle s_2, p_2, o_2 \rangle$  are <u>coalescable</u> if and only if  $\{s_1, o_1\}$  and  $\{s_2, o_2\}$  share at least one common variable.

Intuitively, two triple patterns are coalescable if they have common variables at the subject or object positions. Since a BGP is composed of triple patterns, we can extend coalescability to BGPs, where we require some of their constituent triple patterns to be coalescable.

Definition 4 (Coalescable BGPs): We say that the BGPs  $b_1$ and  $b_2$  are <u>coalescable</u> if there exist  $t_{i_1} \in b_1$  and  $t_{i_2} \in b_2$ such that  $t_1$  and  $t_2$  are coalescable triple patterns. We denote the coalesce operation as +, e.g., the BGP formed from coalescing  $b_1$  and  $b_2$  is denoted as  $b_1 + b_2$ .

*Definition 5 (Basic Graph Pattern (BGP)):* A BGP is recursively defined as follows:

- 1) A triple pattern *t* is a BGP;
- 2) if P<sub>1</sub> and P<sub>2</sub> are coalescable BGPs, P<sub>1</sub> + P<sub>2</sub> is also a BGP. *Definition 6 (Graph Pattern, Group Graph Pattern):*A graph pattern is recursively defined as follows:
- 1) if *P* is a BGP, *P* is a graph pattern;



Fig. 2: An example SPARQL query and its tree expression

- 2) if *P* is a group graph pattern (defined below), *P* is a graph pattern;
- 3) if  $P_1$  and  $P_2$  are both graph patterns,  $P_1$  AND  $P_2$  is also a graph pattern;
- 4) if P<sub>1</sub> and P<sub>2</sub> are both graph patterns, {P<sub>1</sub>} UNION {P<sub>2</sub>}, P<sub>1</sub> OPTIONAL {P<sub>2</sub>} are both graph patterns. Note that {P<sub>i</sub>} denotes a group graph pattern (defined below);

A group graph pattern *P* is recursively defined as follows:

1) If *P* is a graph pattern,  $\{P\}$  is a group graph pattern.

Fig. 2 is an example SPARQL query with six triple patterns  $(t_{1,6})$  and UNION and OPTIONAL operators.

**SPARQL Query—Semantics.** A graph pattern is essentially an expression that contains triple patterns and the left-associative operators AND, UNION and OPTIONAL, which accept graph patterns as their operands. AND and OPTIONAL are binary operators, while UNION can have two or more operands. Such an expression can be equivalently represented by a tree, where each leaf node represents a triple pattern and each internal node represents an operator. The positioning of each node in the tree is determined by the priority of the operators: AND = OPTIONAL  $\prec$  UNION. Curly braces override the standard order of operations and define the scope of the operands. Fig. 2b shows such a tree expression of the query in Fig. 2a.

A graph pattern *P* is matched on an RDF dataset *D* (denoted by  $[[P]]_D$ ) to produce a bag (i.e., multi-set) of mappings  $\{\mu_1, \mu_2, \ldots, \mu_n\}$ , which may contain duplicate mappings. A mapping  $\mu : V \mapsto U$  is a partial function from *V* to  $(I \cup L)$ , where *V* represents the variables that appear in the query, and *I* and *L* denote the sets of IRI and literals, respectively. The set of variables appearing in mapping  $\mu$  is denoted by  $dom(\mu)$ . The two mappings  $\mu_1$  and  $\mu_2$  are defined to be *compatible* (denoted by  $\mu_1 \sim \mu_2$ ) if and only if for all variables  $\nu \in dom(\mu_1) \cap dom(\mu_2)$  satisfying  $\mu_1(\nu) = \mu_2(\nu)$ . Intuitively, this means that the common variables of  $\mu_1$  and  $\mu_2$  are compatible,  $\mu_1 \cup \mu_2$  is also a mapping. If the two mappings  $\mu_1$  and  $\mu_2$  are *mapped* to the case as  $\mu_1 \not\sim \mu_2$ .

We denote two bags of mappings by  $\Omega_1$  and  $\Omega_2$ , and define several operators on bags as follows:

TABLE I: An example RDF dataset

- E	a 1.	- 11	-14	
	Subject	Predicate	Object	
İ	dbr:George_WBush	foaf:name	"George Walker Bush"@en	
[	dbr:George_WBush	rdfs:label	"George W. Bush"@en	
ſ	dbr:George_WBush	dbo:wikiPageWikiLink	dbr:President_of_the_United_States	
ſ	dbr:Bill_Clinton	foaf:name	"Bill Clinton"@en	
ſ	dbr:Bill_Clinton	dbo:wikiPageWikiLink	dbr:President_of_the_United_States	
ſ	dbr:Bill_Clinton	dbp:birthDate	"1946-08-19"^ ^ xsd:date	
[	dbr:Bill_Clinton	owl:sameAs	fbp:Clinton_William_Jefferson_1946-	

1) 
$$\Omega_1 \bowtie \Omega_2 = \{\mu_1 \cup \mu_2 \mid \mu_1 \in \Omega_1 \land \mu_2 \in \Omega_2 \land \mu_1 \sim \mu_2\}.$$

- 2)  $\begin{array}{l} \Omega_1 \cup_{bag} \Omega_2 = \{\mu_1 \mid \mu_1 \in \Omega_1\} \bigcup_{bag} \{\mu_2 \mid \mu_2 \in \Omega_2\}. \\ 3) \quad \Omega_1 \setminus \Omega_2 = \{\mu_1 \in \Omega_1 \mid \forall \mu_2 \in \Omega_2 : \mu_1 \nsim \mu_2\}. \\ 4) \quad \Omega_1 \ \bowtie \Omega_2 = (\Omega_1 \bowtie \Omega_1) \bigcup_{bag} (\Omega_1 \setminus \Omega_1) \end{array}$

Note that the operators above all preserve duplicate elements, as they follow the bag semantics.

We define how to evaluate graph patterns according to W3C's standard semantics for SPAROL as follows.

Definition 7 (Evaluation of Graph Patterns on an RDF dataset): The evaluation of graph patterns P on an RDF dataset D, denoted as  $[P]_D$ , is recursively defined as follows, where the order of the operators follow that in the tree expression:

- 1) If *P* is a triple pattern *t*,  $[[P]]_D = {\mu \mid var(t) = dom(\mu) \land$  $\mu(t) \in D$  (*var*(*t*) represents all variables occurring in t, and  $\mu(t)$  mean that all variables appearing in t are replaced by  $\mu$ ).
- 2) If  $P = \{P_1\}, [[P]]_D = [[P_1]]_D$ .
- 3) If  $P = P_1$  AND  $P_2$ ,  $[[P]]_D = [[P_1]]_D \bowtie [[P_2]]_D$ .
- 4) If  $P = P_1$  UNION  $P_2$ ,  $[[P]]_D = [[P_1]]_D \bigcup_{bag} [[P_2]]_D$ .
- 5) If  $P = P_1$  OPTIONAL  $P_2$ ,  $\llbracket P \rrbracket_D = \llbracket P_1 \rrbracket_D \bowtie \llbracket P_2 \rrbracket_D$ .

#### **IV. PLAN REPRESENTATION: BGP-BASED EVALUATION TREE**

The most straightforward approach for evaluating a graph pattern P is to employ a bottom-up strategy on the tree representation. In each step, we either evaluate a triple pattern, or perform an operator (AND, UNION, or OPTIONAL). This tree-based evaluation strategy strictly follows the SPARQL semantics discussed in Sec. III, but has inherent performance limitations due to the large number of intermediate results generated for each triple pattern at the leaf nodes. To illustrate this, consider the simple SPARQL query in Fig. 3. Note that the outermost group graph pattern of this query only contains a BGP. Following the tree-based method, we first need to obtain  $[[t_1]]_D$  and  $[[t_2]]_D$ . Obviously, the triple pattern  $t_2$  will generate a large number of intermediate results, since most persons in the database have their birth dates as an attribute.

It is more desirable to use BGP evaluation as the basic building block for executing SPARQL queries, employing an optimized BGP query evaluation method such as those used in RDF-3x [11], SW-store [6], gStore [4] and Jena [8]. Thus, in our approach, we design a BGP-based Evaluation Tree (BE-tree) to represent the intermediate form between SPARQL queries and their RA-based query plans.

# A. BE-Tree & RA-Based Plan Structure

The BE-tree is a conceptually simple syntax parse tree of the SPAROL dialect considered in this paper, which is straightforward to construct and intuitive to understand. We use the BE-tree as an intermediate form between a SPARQL query and its RA-based plan.

Definition 8 (BGP-based Evaluation Tree (BE-tree)): Given a group graph pattern Q, its corresponding BE-tree T(Q) is recursively defined as follows:

- The root of T(Q) is a group graph pattern node (Def. 6) representing the query  $\overline{Q}$ ;
- An internal node of T(Q) can be one of {UNION, OPTIONAL, group graph pattern} nodes:
  - A UNION node represents the UNION expression that links two or more group graph patterns, called UNION'ed group graph patterns. It has two or more child nodes, which are all group graph pattern nodes;
  - An OPTIONAL node represents the OPTIONAL expression that links OPTIONAL-left and OPTIONALright graph patterns. It has exactly one child node: the OPTIONAL-right graph pattern, which is a group graph pattern node;
- Each leaf node of T(Q) is a BGP node (Def. 5).

After constructing a BE-tree from parsing a SPARQL query, all coalescable sibling triple patterns and BGPs whose parent is a group graph pattern node are coalesced until no further coalescing can be performed, forming maximal BGPs. When the triple patterns of siblings or BGPs are on different sides of an OPTIONAL node, they can only be coalesced when the OPTIONAL is well-designed [21].

As a concrete example, the BE-tree of the query in Fig. 2a is given in Fig. 4. Note that the triple patterns  $t_1$  and  $t_6$ are coalesced to form a BGP node; no other triple patterns cannot be coalesced, and thus form individual BGP nodes.

The RA-based plan, which dictates how the query is executed, is defined as follows.

Definition 9 (RA-Based Plan): Given a group graph pattern Q, its corresponding RA-based plan Plan(Q) is defined as follows:

- Each internal node of Plan(Q), including the root node, is an AND, OPTIONAL, or UNION node;
- Each leaf node of Plan(Q) is a BGP node (Def. 5).

Compared with the tree expression in Fig. 2b, the RAbased plan has the same types of internal nodes, but has BGPs as leaf nodes instead of triple patterns. For convenience, we denote the internal nodes as the graph pattern



Fig. 3: Inefficiency of binary-tree-based query evaluation



 $\begin{array}{c|c} & & & \\ & & & \\ \hline BGP \\ t_1, t_6 \\ \hline BGP \\ t_2 \\ \hline t_2 \\ \hline BGP \\ t_3 \\ \hline \end{array} \\ \begin{array}{c} BGP \\ BGP \\ t_3 \\ \hline \end{array} \\ \begin{array}{c} BGP \\ BGP \\ t_3 \\ \hline \end{array} \\ \begin{array}{c} BGP \\ t_5 \\ \hline \end{array} \\ \end{array} \\ \begin{array}{c} BGP \\ t_5 \\ \hline \end{array} \\ \end{array}$  \\ \begin{array}{c} BGP \\ t\_5 \\ \hline \end{array} \\ \end{array} \\ \begin{array}{c} BGP \\ t\_5 \\ \hline \end{array} \\ \end{array} \\ \begin{array}{c} BGP \\ t\_5 \\ \hline \end{array} \\ \end{array} \\ \end{array}

Fig. 4: Example BE-tree.

Fig. 5: Example RA-based plan.

operators' respective RA operators, i.e., AND as  $\bowtie$ , UNION as  $\bigcup_{bag}$ , and OPTIONAL as  $\bowtie$ . For example, Fig. 5 shows the RA-based plan converted from the BE-tree Fig. 4. To evaluate a SPARQL query, its RA-based plan is traversed in a depth-first manner, during which the BGPs are evaluated and their results are combined according to the operators' semantics, as shown in Alg. 1.

Algorithm 1: BGP-based query evaluation					
	Input: RDF dataset D, RA-based plan Plan(Q)				
	<b>Output:</b> [[Q]] <sub>D</sub>				
1	BGPBasedEvaluation(D, Plan(Q).root);				
2	Function BGPBasedEvaluation (D, v):				
3	if v is of type $\bowtie$ then				
4	$r \leftarrow \text{BGPBasedEvaluation}(D, v.child[0]);$				
5	$r \leftarrow r \bowtie BGPBasedEvaluation(D, v.child[1]);$				
6	return r;				
7	else if v is of type $\bigcup_{hag}$ then				
8	$r \leftarrow \text{BGPBasedEvaluation}(D, v.child[0]);$				
9	for i in $1, 2, \cdots$ , num(child) do				
10	$r \leftarrow r \bigcup_{bag} BGPBasedEvaluation(D, v.child[i]);$				
11	return r;				
12	else if v is of type $\exists \forall$ then				
13	$r \leftarrow \text{BGPBasedEvaluation}(D, v.child[0]);$				
14	$r \leftarrow r \square BGPBasedEvaluation(D, v.child[1]);$				
15	return r;				
16	else if $v$ is a BGP then				
17	return EvaluateBGP $(D, e_i)$				

It is straightforward to convert from a BE-tree to an RAbased plan. Basically, we perform a depth-first search on the BE-tree, prioritizing each node's rightmost children. The RAbased plan is grown from right to left as the BE-tree nodes are visited. A group graph pattern node in a BE-tree will be mapped to one or more  $\bowtie$  nodes in an RA-based plan.

## B. BE-Tree & RA-Based Plan Transformations

In the previous subsection, we invoke the BGP-based evaluation procedure (Alg. 1) on the RA-based plan directly

constructed from BE-tree of the query. However, it is possible to improve the efficiency of query evaluation by altering the plan. We achieve this by making certain semanticspreserving *transformations*.

1) Goals: Our aim is to transform the original RA-based plan so that the resulting RA-based plan has the following properties:

- *Validity:* the resulting RA-based plan should maintain the previously defined tree structure and have the same node types.
- *Efficiency:* the evaluation of the resulting RA-based plan should be more efficient than the original RA-based plan.

2) Semantics-Preserving Transformations: In order to optimize for query execution efficiency while maintaining correctness, we need to leverage the inherent semantic equivalences regarding the UNION and OPTIONAL, according to the following theorems backed by RA. We first introduce these transformations on the BE-tree, then map them to RA-based plans.

*Theorem 1:* For any graph pattern  $P_1$ ,  $P_2$ ,  $P_3$  and any RDF dataset D, we have

 $\llbracket P_1 \text{ AND } (P_2 \text{ UNION } P_3) \rrbracket_D = \llbracket (P_1 \text{ AND } P_2) \text{ UNION } (P_1 \text{ AND } P_3) \rrbracket_D.$ 

Proof 1:

 $\begin{bmatrix} P_1 \text{ AND } (P_2 \text{ UNION } P_3) \end{bmatrix}_D$   $= \begin{bmatrix} P_1 \end{bmatrix}_D \bowtie \begin{bmatrix} P_2 \text{ UNION } P_3 \end{bmatrix}_D$   $= \begin{bmatrix} P_1 \end{bmatrix}_D \bowtie (\begin{bmatrix} P_2 \end{bmatrix}_D \cup_{bag} \begin{bmatrix} P_3 \end{bmatrix}_D)$   $= (\begin{bmatrix} P_1 \end{bmatrix}_D \bowtie \begin{bmatrix} P_2 \end{bmatrix}_D) \cup_{bag} (\begin{bmatrix} P_1 \end{bmatrix}_D \bowtie \begin{bmatrix} P_3 \end{bmatrix}_D)$   $= \begin{bmatrix} P_1 \text{ AND } P_2 \end{bmatrix}_D \cup_{bag} \begin{bmatrix} P_1 \text{ AND } P_3 \end{bmatrix}_D$   $= \begin{bmatrix} (P_1 \text{ AND } P_2) \text{ UNION } (P_1 \text{ AND } P_3) \end{bmatrix}_D$ 

Note that Thm. 1 is also trivially extendable to UNION nodes with more than two child nodes.

*Theorem 2:* For any graph pattern  $P_1$ ,  $P_2$  and any RDF dataset *D*, we have

$$\llbracket P_1 \text{ OPTIONAL } P_2 \rrbracket_D = \llbracket P_1 \text{ OPTIONAL } (P_1 \text{ AND } P_2) \rrbracket_D.$$

Proof 2:

$$\begin{split} & [\![P_1 \text{ OPTIONAL } (P_1 \text{ AND } P_2)]\!]_D \\ &= ([\![P_1]\!]_D \bowtie [\![P_1 \text{ AND } P_2]\!]_D) \cup_{bag} ([\![P_1]\!]_D \setminus [\![P_1 \text{ AND } P_2]\!]_D) \\ &= ([\![P_1]\!]_D \bowtie ([\![P_1]\!]_D \bowtie [\![P_2]\!]_D)) \\ & \cup_{bag} ([\![P_1]\!]_D \setminus ([\![P_1]\!]_D \bowtie [\![P_2]\!]_D)) \\ &= ([\![P_1]\!]_D \bowtie [\![P_2]\!]_D) \cup_{bag} ([\![P_1]\!]_D \setminus [\![P_2]\!]_D) \\ &= [[\![P_1 \text{ OPTIONAL } P_2]\!]_D. \end{split}$$

These two equivalences correspond to two semanticspreserving transformations on the BE-tree: that of *merging* a node with the child nodes of its sibling UNION node, and that of *injecting* a node into the child node of its sibling OPTIONAL node, defined as follows.

Definition 10 (Merge transformation): A merge transformation is performed on a graph pattern  $P_1$ 's node and one of its sibling UNION nodes, the child nodes of which represents the group graph patterns  $P_2, P_3, \dots, P_n$ , when

- 1)  $P_1$  is a BGP node;
- 2) At least one of the group graph patterns in  $P_2, P_3, \dots, P_n$  has a BGP child node that is coalescable with  $P_1$ .

The merge transformation consists of the following steps:

- 1) Insert  $P_1$  as the leftmost child node of  $P_2, P_3, \dots, P_n$ ;
- Coalesce P<sub>1</sub> with the other BGP child nodes if possible, until all the BGP nodes are maximal;
- 3) Remove  $P_1$  from its original position.

Definition 11 (Inject transformation): An inject transformation is performed on a graph pattern  $P_1$ 's node and one of its sibling OPTIONAL nodes to its right, the child node of which represents the group graph pattern  $P_2$ , when

- 1)  $P_1$  is a BGP node;
- 2)  $P_2$  has a BGP child node that is coalescable with  $P_1$ . The inject transformation consists of the following steps:
- 1) Insert  $P_1$  as the leftmost child node of  $P_2$ ;
- 2) Coalesce  $P_1$  with the other BGP child nodes if possible, until all the BGP child nodes are maximal.



Fig. 6: Example *Inject* Transformation.

Fig. 7: Example *Merge* Transformation.

*Example.* In Fig. 6 and 7, we give examples of these transformations' effects on SPARQL queries targeting DBpedia.

In Fig. 6,  $b_4$  is a grandchild BGP node of the OPTIONAL node to the right of  $b_1$ , and  $b_1$  and  $b_4$  are coalescable. Therefore, the available inject transformation will coalesce  $b_4$  with  $b_1$ , which can help improve efficiency. According to the original BE-tree,  $b_4$  is directly evaluated, and the results are left-outer-joined with those of  $b_1$ . Since a large number of entities have the ?sameAs relation, which denotes the equivalence between references to the same real-world object, b<sub>4</sub> has many matches, causing both its evaluation and the left-outer-join to be costly. However, presidents of the United States is a minority of the entities, making  $b_1$  highly selective. After the *inject*, we can rely on the underlying evaluation engine to efficiently evaluate the coalesced  $b_1b_4$  by choosing a join order that evaluates the much more selective  $b_1$  first. The left-outer-join is also rendered less expensive due to the decrease in the number of results of  $b_1b_4$  compared with  $b_4$ .

This example also explains why the transformations are only defined on coalescable BGPs. If no coalescing happens, the repeated evaluation of the merged or injected BGP will incur overhead.

However, not all available transformations can help improve efficiency, which necessitates enumeration and costbased selection. Fig. 7 shows an available *merge* transformation on an example UNION query, which merges the BGP  $b_1$  with its sibling UNION node. Since  $b_1$  has low selectivity, merging it does not accelerate BGP evaluation or reduce the number of intermediate results, and even incurs extra overhead because it now has to be evaluated twice.

When mapped onto the RA-based plan, the scope of a transformation is not as obvious as sibling nodes in the BE-tree. To search for UNIONS and OPTIONALS viable for transforming with a BGP, we need to search the BGP's ancestors until finding the  $\bowtie$  operator closest to the root without passing through any  $\bigcup_{bag}$  node. All the  $\bigcup_{bag}$ 's and  $\bowtie$ 's reachable from this  $\bowtie$  through only  $\bowtie$ 's are exactly this BGP's siblings in the BE-tree and can be considered for transformation. We denote the set of  $\bigcup_{bag}$ 's and  $\bowtie$ 's that can be considered by a BGP *b* as scope(*b*).

## V. COST-DRIVEN PLAN SELECTION

In this section, we introduce the cost-driven approach selecting the expectedly most efficient transformations to be performed on an RA-based plan.

Our cost-driven approach operates on a higher level than BGP evaluation, but still relies on estimations of the evaluation costs and result sizes of BGPs, which are obtainable from the underlying BGP evaluation engine [22].

## A. Cost Models

1) Cost Model for SPARQL-UO: SPARQL-UO query execution cost is made up of two main components: the cost of evaluating BGPs and of computing graph pattern operators:  $\bowtie, \bigcup_{bag}$ , and  $\bowtie$ . We are primarily concerned with the cost difference caused by a transformation, which we call  $\Delta$ -cost. A transformation is expected to improve efficiency only

when its  $\Delta$ -cost is negative. In the following, we discuss how to estimate the  $\Delta$ -cost.



Fig. 8: Merge & Inject Transformations on an RA-Based Plan.

a) Merge transformation: Consider the RA-based plan in Fig. 8a. The merge transformation shown as the red arrow will cause the following structural changes in the RA-based plan: the outer BGP will be coalesced with the  $\bigcup_{bag}$ 's child BGPs, and if some children of the  $\bigcup_{bag}$  are not BGPs or are not coalescable with the outer BGP, new  $\bowtie$ 's will join them with the outer BGP. The transformation will also affect the cost of the  $\bigcup_{bag}$ , since the operands' result sizes will change. The cost of all the ancestor operators of the  $\bigcup_{bag}$  will also change if the result size of the  $\bigcup_{bag}$  changes, but we do not account for them for efficiency. When estimating the  $\Delta$ -cost of a merge operation, we only need to account for the cost of the affected parts in the RA-based plan before and after the operation.

Suppose a merge transformation concerns the outer BGP b and the  $\bigcup_{bag}$  node u; the set of child nodes of u in the RA-based plan is denoted as child(u). We can express the aforementioned affected parts' cost before and after the merge by the following formulas:

$$cost_{BGP}(b,u) = cost(b) + \sum_{b_i \in child(u)} cost(b_i)$$
(1)

$$cost_{algebra}(b,u) = cost(u)$$
 (2)

$$cost_{BGP}(merge(b, u)) = \sum_{b_i \in child(u)} cost(b + b_i)$$
 (3)

$$cost_{algebra}(merge(b, u)) = cost(u') + \sum_{p_i \in child(u)} cost(\bowtie(b, p_i))$$
(4)

$$\Delta cost(b, u) = cost_{BGP}(b, u) + cost_{algebra}(b, u)$$

$$- cost_{BGP}(merge(b, u))$$

$$- cost_{algebra}(merge(b, u))$$
(5)

where Eq. 1 and 3 represent the cost of the affected BGPs, i.e., the outer BGP b and the child BGPs of u that are coalescable with b, denoted as  $b_i$ , before and after the merge operation, since b is coalesced with each  $b_i$  during the merge; while Eq. 2 and 4 represent the cost of the affected algebraic operators, including u and the newly introduced  $\bowtie'$ s. Thus, the  $\Delta$ -cost of this merge operation

is the difference between the summed cost of the affected BGPs and algebraic operators before and after the operation.

The cost of the BGPs can be directly obtained from the BGP evaluation engine, while the cost of the algebraic operators are functions on the result sizes of their operands. These functions may differ with physical implementations. In our experiments, to fit the system we choose for implementation, the cost of  $\bowtie$  is set as the product of its arguments' result sizes, and the cost of the  $\bigcup_{bag}$  is set as the sum of its arguments' results sizes. A BGP node's result size estimate can be obtained from the BGP evaluation engine. The result sizes of other types of nodes need to be estimated based on an assumed distribution of data. In our experiments, we estimate the result size of any  $\bowtie$  or  $\bowtie$  as the product of their operands' result sizes, and the result sizes of  $\bigcup_{bag}$  to be the sum of its operands' result sizes.

b) Inject transformation: An inject transformation such as the red arrow in Fig. 8b will cause a BGP in the subtree rooted at the right child of the  $\bowtie$  node to coalesce with the outer BGP. In practice, we choose the first coalescable BGP as the target for injection for convenience. The transformation will also affect the cost of the  $\bowtie$  as well as all the  $\bowtie$ 's on the path from the  $\bowtie$  to the coalesced BGP.

Suppose an inject transformation concerns the outer BGP b and the  $\bowtie$  node o due to the change in result sizes. The following formulas show the affected parts' cost before and after injecting:

$$cost_{BGP}(b,o) = cost(b_i)$$
(6)

$$cost_{algebra}(b,o) = cost(o) + \sum_{a_i \in p(o,b_i)} cost(a_j)$$
(7)

$$cost_{BGP}(inject(b, o)) = cost(b + b_i)$$
 (8)

$$cost_{algebra}(inject(b,o)) = cost(o') + \sum_{a'_i \in p(o,b_i)} cost(a_j')$$
(9)

$$\Delta cost(b, o) = cost_{BGP}(b, o) + cost_{algebra}(b, o)$$
(10)  
$$- cost_{BGP}(inject(b, o))$$
  
$$- cost_{algebra}(inject(b, o))$$

where  $b_i$  in Eq. 6 denotes the BGP in *o*'s right child's subtree; after injecting, it coalesces with *b* as shown in Eq. 8.  $a_j$  denotes the  $\bowtie$ 's on the path  $p(o, b_i)$  from the  $\bowtie$  to the coalesced BGP. o' and  $a_j'$  denotes the  $\bowtie$  and  $\bowtie$ 's whose cost changes due to the change in cardinality.

The cost and cardinality of the BGPs and the algebraic operators involved are estimated in the same way as for merge transformations.

2) Cost Model for BGPs: We briefly introduce the BGP cost models employed by gStore, one of the systems we select for implementation, for completeness. The evaluation of BGPs consists of joins. Thus the cost of a BGP plan T is the sum of each join *j*'s cost:

$$cost(T) = \sum_{j \in T} cost(j)$$

BGP evaluation in gStore uses the worst-case optimal (WCO) join. For each result tuple on the existing vertices, all such edges need to be scanned at least once to check whether this tuple can be extended to match the newly extended vertex. Suppose the set of existing vertices is  $\{v_1, \dots, v_{k-1}\}$ , and the newly extended vertex is  $v_k$ . The cost of a WCO join can then be estimated as follows:

$$cost(WCOJoin(\{v_1, \dots, v_{k-1}\}, v_k))) = card(\{v_1, \dots, v_{k-1}\}) \times \min_{i \in [1, k-1]} average\_size(v_i, p)$$

where  $card(\{v_1, \dots, v_{k-1}\})$  indicates the estimated cardinality, and average\_size( $v_i, p$ ) indicates the average number of edges (*i.e.*, triples) with p as predicate and  $v_i$  as subject or object.

The cardinality estimation of query vertex sets starts from single triple patterns, whose exact cardinality can be obtained from the RDF store. Each time a new query vertex is added to the set, we sample the candidate result set, and collate how many result tuples can be generated from the sample by extending to the new query vertex. The estimated cardinality is updated by scaling up the previous estimate:

$$card(V_k) = \max(\frac{\#extend}{\#sample} \times card(V_{k-1}), 1)$$

#### B. Cost-Driven Transformation

In this subsection, we discuss RA-based plan transformation selection algorithms that leverage the cost model discussed above to decide on transformations for obtaining the most efficient query plan for execution.

1) Transforming a Single BGP: We first concentrate on the simpler case where only transformations concerning a BGP are considered, in which case the algorithm for transformation selection is shown in Alg. 2. The algorithm first finds the sets of  $\bigcup_{hag}$ 's and  $\bowtie$ 's in the RA-based plan that can perform a transformation with the given BGP (Line 2), the details of which will be introduced below. According to Thm. 1 and 2, a merged BGP is removed from its original position, while an injected BGP maintains its original occurrence. This means that a BGP can only be merged with one of the  $\bigcup_{bag}$ 's, but can be injected into multiple 3%. Therefore, in order to decide on a merge transformation, we need to look at all the viable  $\bigcup_{bag}$ 's and choose the transformation that incurs the lowest  $\Delta$ -cost (Lines 3-11 and 14-17). On the other hand, inject transformations are mutually independent, so we scan over each  $\mathbb{N}$  and decide individually which ones should be transformed based on the  $\Delta$ -cost (Lines 12-13 and 18-22).

We further explain how to find the sets of  $\bigcup_{bag}$ 's and M's viable for transformation with a given BGP, i.e., the FindRelevantUO subroutine in Alg. 2. Assuming welldesigned OPTIONALS [21] throughout the query, the set of viable M's consists exactly of those joined with the given BGP via any number of M's or M's, while the set of viable  $\bigcup_{bag}$ 's consists exactly of those joined with the BGP via any number of M's. Therefore, we search upwards for a root of a subtree in the RA-based plan via only M's or only M's and M's (Lines 26-32) and put all the  $\bigcup_{bag}$ 's or M's reachable via these operators from the root into the viable sets (Lines 33-34 and 36-43).

2) Handling an RA-Based Plan: Handling the entire RA-based plan, which often consists of multiple

#### Algorithm 2: Single-BGP Transformation Selection

```
Input: RDF dataset D, RA-based plan P, a BGP b
 1 Function SingleBGPTransform (D,P,b):
         U, O \leftarrow FindRelevantUO(b, P)
 2
         minUnionCost \leftarrow 0
 3
         targetUNION \leftarrow empty node
 4
         foreach u \in U do
 5
 6
               minUnionCostCur \leftarrow DecideMerge (b, u)
               if minUnionCostCur < minUnionCost then
 7
                    minUnionCost ← minUnionCostCur
 8
 9
                    targetUNION \leftarrow u
10
         if minUnionCost < 0 then
11
             Perform merge on subBGPglobal and targetUNION
12
         foreach o \in O do
13
             DecideInject (b, o)
14 Function DecideMerge(b,u):
15
         if constraints are violated then
16
              return 0
         return \triangle cost(b, u) (Eq. 5)
17
18 Function DecideInject(b,o):
         if constraints are violated then
19
20
              return
21
         if \Delta cost(b, o) < 0 (Eq. 10) then
22
              Perform inject on b and o
23
   Function FindRelevantUO(b, P):
24
         U \leftarrow \emptyset, O \leftarrow \emptyset, \text{ curNode } \leftarrow b
25
         localRootU \leftarrow empty node, localRootO \leftarrow empty node
         while curNode is not P's root do
26
               if curNode.parent is \bigcup_{\textit{bag}} or \bowtie and localRootU is an empty node
27
                then
                   localRootU ← curNode
28
               if curNode.parent is \bigcup_{bag} then
29
                    localRootO \leftarrow curNode
30
31
                    break
              curNode ← curNode.parent
32
         SearchSubtree (localRootU, P, U)
33
         SearchSubtree (localRootO, P, O)
34
35
         return U,O
36 Function SearchSubtree (v, P, S):
         if v is an empty node then
37
              v \leftarrow P.root
38
         if The elements in S are of the same type as v then
39
40
              S \leftarrow S \cup \{v\}
41
         else if v is \bowtie or those in S are \bowtie and v is \bowtie then
               SearchSubtree(v.leftChild, P, S)
42
               SearchSubtree (v.rightChild, P, S)
```

BGPs, is challenging because of the possible interdependence between transformations across BGPs. For example, if we consider transforming the group graph pattern  $\{P_1 \text{ OPTIONAL } \{P_2 \text{ OPTIONAL } P_3\}\}$   $\{P_1, P_2 \text{ and } P_3 \text{ are}$ all coalescable BGPs), there are  $2^3$  possible transformations involving whether  $P_1$  is injected into  $P_2$ , whether  $P_2$  is injected into  $P_3$ , and whether  $P_1$  is injected into  $P_3$ . This results in a plan space that is exponential in terms of the depth of the RA-based plan.

In order to balance the time complexity and the efficiency of the transformed plan, we propose a greedy strategy to decide on the transformations on the entire RA-based plan (Alg. 3). Specifically, the algorithm first conducts a topological sort over the RA-based plan (Line 1). The BGPs are then considered in a reverse topological order for transformations (Lines 3-5). In this way, we ensure that the BGPs that are lower in the plan have been appropriately transformed before considering transforming the BGPs higher up in the plan, preventing expensive backtracking.

## Algorithm 3: RA-Based Plan Transformation Selection

	Input: RDF dataset D, RA-based plan P				
1	<pre>Function RAPlanTransform(D, P):</pre>				
2	Do a topological sort on P's nodes				
3	Let B be the BGPs in P sorted in descending topological order				
4	foreach $b \in B$ do				
5	SingleBGPTransform(D, P, b)				

#### VI. QUERY-TIME OPTIMIZATION: CANDIDATE PRUNING

In the previous section, we introduced how to select transformations on the RA-based plan based on cost estimations prior to execution. In this section, we present *candidate pruning*, a query-time optimization incorporated into Alg. 1 to enhance efficiency.



Fig. 9: Candidate pruning for OPTIONAL.

The basic idea of candidate pruning is also drawn from Thm. 1 and 2. The equivalence between the evaluation results implies that the results of the  $\bigcup_{bag}$ 's and  $\mathbb{H}$ 's are constrained by those of the outer graph pattern in term of the common variables. Therefore, when a  $\bigcup_{bag}$  or  $\mathbb{H}$ is encountered during evaluation, we can set the current results on the common variables as *candidate results* when executing the child BGPs of that node. For ease of explication, we show the mechanism of candidate pruning for an OPTIONAL query in Fig. 9, where the results of the variable ?x from the already evaluated graph patterns serve as the candidate results of ?x for the child BGP of the group graph pattern below the OPTIONAL, pruning redundant matchings of ?x that will be materialized if the BGP is evaluated independently.

To achieve a pruning effect, we need to ensure that the size of the candidate results is smaller than the size of the actual results of the BGP. A smaller candidate result size also reduces the overhead incurred by scanning them and setting them as candidates. Therefore, we adopt an adaptive threshold on the candidate result size. The cost model for BGPs (Sec. V-A2) invoked as part of the transformation selection procedure (Alg. 2) provides an estimate of the actual BGP result size, which we employ as the threshold on candidate result size whenever possible. When no such estimate is available, we set the threshold based on the dataset size. (Please refer to Sec. VII for the threshold setting in our experiments.)

To implement candidate pruning, we modify Alg. 1 as follows (Note that the results can be passed as arguments in the form of pointers to prevent expensive copying):

- Add a third argument *cand*, which denotes the candidate results, to the BGPBasedEvaluation function;
- Pass the current results r as the third argument to BGPBasedEvaluation when processing a  $\bigcup_{bag}$  or  $\bowtie$ (Lines 10 and 14);
- Pass *cand* as the third argument to EvaluateBGP (Line 17). Only when the size of *cand* is smaller than the threshold is it set as the candidate results of the BGP.

Tree transformations, which happens before query execution, and candidate pruning, which take effect during execution, are complementary to each other. Prior to execution, high-selectivity BGPs are targeted by merge or inject transformations, which breaks up graph patterns with large overall results that originally cannot be handled by candidate pruning. Meanwhile, while tree transformations are constrained to be performed one BGP at a time due to the vast plan space, candidate pruning can transmit the pruning effect of small results across levels during execution. For example, when processing a query with the group graph pattern { $P_1$  OPTIONAL { $P_2$  OPTIONAL  $P_3$ }},  $P_1$  cannot be injected into  $P_3$  by the greedy transformation strategy even if it is selective, but its results can serve as candidates for  $P_3$ via  $P_2$ . In the special case where there is only a BGP viable for transforming with a UNION or OPTIONAL, performing transformations on them is equivalent to candidate pruning. In this case, tree transformation is skipped to evade the additional overhead.

#### VII. EXPERIMENTS

To evaluate the effectiveness of our approach, we employ the BGP query engines of Jena<sup>1</sup>, Blazegraph<sup>2</sup>, and gStore to implement our BGP-based cost-aware SPARQL-UO evaluation strategy. All the experiments run on Jena have enabled the statistics-based optimizations. We forked a branch from the main branch of gStore and implement our proposed SPARQL-UO optimizer based on it<sup>3</sup>. Experiments are conducted on both synthetic (LUBM [23]) and real (DBpedia<sup>4</sup> [24] and YAGO [25]) RDF datasets, the statistics of which are listed in Table II. Our implementation and all the queries used in our experiments can be found in our GitHub repository<sup>5</sup>. We conduct experiments on a Linux server with an Intel Xeon Gold 6126 CPU @ 2.60GHz CPU and 256GB memory.

#### A. Verification of Optimizations

In this section, we verify the effectiveness of the proposed optimizations in Section IV-B and evaluate the following four approaches:

<sup>&</sup>lt;sup>1</sup>https://github.com/apache/jena.

<sup>&</sup>lt;sup>2</sup>https://github.com/blazegraph/database.

<sup>&</sup>lt;sup>3</sup>Our implementation is available at https://github.com/SoftlySpoken/gStore-UO-opt.

<sup>&</sup>lt;sup>4</sup>The DBpedia data dump that we use is V3.9, which is downloadable at http://downloads.dbpedia.org/3.9/en/. We use the concatenation of all the N-Triples files.

<sup>&</sup>lt;sup>5</sup>https://github.com/SoftlySpoken/gStore-UO-opt.

TABLE II: Datasets Statistics

Datasets	triples	entities	predicates	literals
LUBM	534,355,247	86,990,882	18	44,658,530
DBpedia	830,030,460	96,375,582	57,471	59,825,935
YAGO	230,677,128	99,313,763	108	153,988,292

TABLE III: Query Statistics on LUBM

		Query	Туре	Count <sub>BGP</sub>	Depth	$ \llbracket Q \rrbracket_D $
		q1.1	U	9	2	645,666
		q1.2	0	3	2	44,653,510
	Group 1	q1.3	0	4	4	76
		q1.4	0	4	4	5,583
		q1.5	UO	6	3	4,348
		q1.6	UO	9	3	37
1		q2.1	0	3	1	4,176,432
	Group 2	q2.2	0	4	3	8,698
		q2.3	0	4	3	13,124,940
		q2.4	0	2	3	10
		q2.5	0	2	2	10
		q2.6	Ó	2	2	7

- The baseline (abbreviated as base), i.e., the base versions of the systems, which invoke the BGP-based query evaluation method (Alg. 1);
- 2) Tree transformation (abbreviated as TT), which transforms the original RA-based plan by Alg. 3 and then invokes Alg. 1 on it;
- Candidate pruning (abbreviated as CP), which invokes Alg. 1 augmented with candidate pruning (Sec. VI) on the original RA-based plan, using a fixed threshold of 1% of the total number of triples in the database;
- 4) The full version that coordinates tree transformation and candidate pruning (abbreviated as full), using an adaptive threshold on the candidate result size.

Since there is no benchmark tailored for SPARQL-UO queries to our knowledge, we construct a mini-benchmark with realistic semantics and varying complexities, containing six queries on LUBM, DBpedia, and YAGO, respectively, denoted as q1.1-1.6 in the following and given in Appendix A of [26]. Let Q be the outermost group graph pattern in the query. To measure the complexity of a query, we define two metrics: (1) the BGP count ( $Count_{BGP}(Q)$ ), and (2) the maximum depth of nested group graph patterns (Depth(Q)).

 $Count_{BGP}(P)$  of a graph pattern P is recursively defined:

1) If P is a BGP, 
$$Count_{BGP}(P) = 1$$
.

- 2) If  $P = \{P_1\}$ ,  $Count_{BGP}(P) = Count_{BGP}(P_1)$ .
- 3) If  $P = P_1$  AND  $P_2$  or  $P_1$  UNION  $P_2$  or  $P_1$  OPTIONAL  $P_2$ ,  $Count_{BGP}(P) = Count_{BGP}(P_1) + Count_{BGP}(P_2)$ .

Depth(P) of a graph pattern P is recursively defined as follows:

- 1) If P is a BGP, Depth(P) = 0.
- 2) If  $P = \{P_1\}$ ,  $Depth(P) = Depth(P_1) + 1$ .
- 3) If  $P = P_1$  AND  $P_2$  or  $P_1$  UNION  $P_2$  or  $P_1$  OPTIONAL  $P_2$ ,  $Depth(P) = max(Depth(P_1), Depth(P_2)).$

Suppose *P* is the outermost group graph pattern of query *Q*, we have  $Count_{BGP}(Q) = Count_{BGP}(P)$ , Depth(Q) = Depth(P). Group 1 in Tables III and IV summarizes the

TABLE IV: Query Statistics on DBpedia

	Query	Туре	Count <sub>BGP</sub>	Depth	$ \llbracket Q \rrbracket_D $
	q1.1	U	6	2	153,325
	q1.2	UO	4	3	610,434
Group 1	q1.3	0	5	5	1,192
	q1.4	UO	7	5	92,041
	q1.5	UO	6	3	3,699,995
	q1.6	UO	10	4	176
	q2.1	0	5	3	490,876
	q2.2	0	2	2	55,054
Group 2	q2.3	0	2	2	61,318
Gloup 2	q2.4	0	3	2	4,757
	q2.5	Ó	2	2	5,330
	q2.6	0	9	2	36

statistics and the result sizes of the queries used in this subsection.

We report the query execution time and the time spent carrying out the tree transformations for TT and full. The performance of our approaches on LUBM, DBpedia, and YAGO is shown in Fig. 10. The absence of a bar indicates an out-of-memory or timeout error on the query. A query is considered timed-out if the execution time exceeds  $2 \times 10^6$  milliseconds.

The trends of the results across gStore, Jena, and Blazegraph are similar, showing the adaptability of our approach regardless of the underlying BGP execution engine. Both of our proposed optimizations are shown to be effective since TT, CP and full perform better than base on all queries. TT and CP can be more advantageous on different queries and datasets than the other. Their optimization effects are cumulative when combined: full performs best all queries and datasets (except on q1.2 on gStore where CP beats full by a small margin, and on q1.3 and q1.6 on Blazegraph, where our optimizations' performance gains just do not compensate for their extra overhead), beating the baseline by up to over an order of magnitude. Our optimized approaches also consume less memory. While base runs out of memory on 13 out of 24 queries, full successfully runs all the queries.

In the following, we try to draw some conclusions about the applicability of our optimizations to different SPARQL-UO queries.

When **TT** is effective. q1.1 on DBpedia (Listing 15, Appendix A of [26]) is a query on which TT is effective, but CP is not. In this query, two UNION clauses are given first (Lines 2-3), whose child BGPs all have low selectivity. There is no high-selectivity graph pattern before them to enable CP. However, TT can merge the high-selectivity BGP in Lines 5-8 with the UNION clause in Line 3 to accelerate query processing and reduce memory overhead, as evidenced in Fig. 10. q1.2 on LUBM and q1.2 on DBpedia



Fig. 10: Verification of optimizations.

also belong to this category. (Note that q1.2 on LUBM corresponds to the special case mentioned in Section VI, where there is only a BGP before an OPTIONAL clause, thus TT and CP have a similar effect.)

When **CP** is effective. q1.3 on LUBM (Listing 4, Appendix A of [26]) is a query on which CP is effective, but TT is not. In

this query, the BGP in Line 2 has high selectivity, followed by nested OPTIONALs with low-selectivity child BGPs. TT can inject the BGP into the outermost OPTIONAL but cannot reach the inner OPTIONALs, thus having limited effect. However, CP can carry the small number of results into the innermost OPTIONAL and set them as candidates to accelerate query processing. q1.3-4 on LUBM and q1.3-4 on DBpedia also belong to this category.

When **TT** and **CP** are jointly effective. q1.6 on LUBM (Listing 7, Appendix A of [26]) is a query on which TT and CP work complementarily, causing full to perform much better than TT and CP. In this query, the BGP in Lines 2-3 has high selectivity, while the BGP in Line 4 has relatively low selectivity. Upon obtaining their considerably large results, CP has limited effect on the following UNION clauses. TT, however, can pick the high-selectivity BGP to merge with the UNION in Line 5. Having executed the graph patterns up to Line 6, CP can accelerate the processing of upcoming OPTIONALs. q1.1 and q1.5 on LUBM and q1.5 and q1.6 DBpedia also belong to this category.

For a quantitative perspective on the optimization effects, we define the join space of a graph pattern JS(P) as follows:

- 1) If *P* is a BGP,  $JS(P) = |\llbracket P \rrbracket_D|$ .
- 2) If  $P = \{P_1\}, JS(P) = JS(P_1).$
- 3) If  $P = P_1$  AND  $P_2$  or  $P_1$  OPTIONAL  $P_2$ ,  $JS(P) = JS(P_1) \times JS(P_2)$ .
- 4) If  $P = P_1$  UNION  $P_2$ ,  $JS(P) = JS(P_1) + JS(P_2)$ .

The join space of a query estimates the largest intermediate result size that is materialized during the execution of this query. Therefore, it is indicative of both the query's execution time and memory overhead. We plot the execution time of all the queries on gStore and Jena (the yaxis on the left) with their respective join spaces (the yaxis on the right) in Fig. 11. Across the tested approaches, these three metrics show a similar trend. On all the queries, the join spaces of TT and CP are smaller than those of base, and full has the smallest join space overall, which corroborates the qualitative analysis above.

#### B. Comparison with State-of-the-Art

The only work that considers SPARQL with OPTIONAL query optimization is LBR [3], with which we compare our full approach. We implement LBR in C++ and experiment on the queries provided in LBR [3] on LUBM and DBpedia, listed as q2.1-2.6 (Appendix A, [26]). The statistics of these queries are given in the second group in Tables III and IV. q2.1-2.3 are complex with multiple nested group graph patterns, each containing a low-selectivity BGP followed by an OPTIONAL with a single low-selectivity child BGP. Meanwhile, q2.4-2.6 are simple without nested group graph patterns, and their outermost group graph pattern has a high-selectivity BGP followed by an OPTIONAL.

The total response time of full and LBR are shown in Fig. 13. full is significantly faster than LBR on all queries,



Fig. 12: Query execution time (ms) of full on LUBM datasets of different sizes ("B" is short for billion).



Fig. 13: Comparison with state-of-the-art on LUBM.

and the improvement on q2.4-2.6 is more significant than on q2.1-2.3. This is because candidate pruning can take advantage of the high-selectivity BGPs in q2.4-2.6, while q2.1-2.3 does not contain high-selectivity BGPs. (Note that since all the group graph patterns in q2.1-2.6 contain a BGP followed by an OPTIONAL clause, they correspond to the special case mentioned in Sec. VI where tree transformation and candidate pruning are equivalent, hence only candidate pruning is performed.) The results show that when candidate pruning takes effect, it is more efficient than LBR's heavy-weight pruning strategies. On q2.1-2.3, full is still faster than LBR since its BGP-based evaluation scheme is more efficient than LBR's separate treatment of triple patterns.

In summary, our approach outperforms LBR on OPTIONAL queries, despite LBR being specially optimized for OPTIONAL.

## C. Scalability Study

Lastly, we evaluate how well our approach scales to larger datasets. By setting the scaling factor of LUBM, i.e., the number of universities, we generate three more LUBM datasets with 1, 1.5 and 2 billion triples, respectively. We run the full approach on q1.1-q1.6 on these datasets and plot how the execution time changes with the dataset size on each query in Fig. 12.

Our approach scales almost linearly to the number of triples in the datasets. The growth rate of the query execution time correlates with each query's result sizes: the execution time of queries with larger result sizes grows faster with the dataset size. (The result sizes of q1.3-1.6 on larger LUBM datasets are equal to those shown in Tab. III, while those of q1.1-1.2 grow linearly.)

#### VIII. CONCLUSION

The proliferation of knowledge graph applications has generated increasing RDF data management problems. In this paper, we focus on how to optimize SPARQL queries with UNION and OPTIONAL clauses (SPARQL-UO for short). Making use of existing BGP query evaluation modules in SPARQL engines, we propose a series of costdriven transformations on the BGP-based evaluation tree (BE-tree). These optimizations can significantly reduce the search space and intermediate result sizes, and thus improve both the time and space efficiency of SPARQL-UO query evaluation. We experimentally validate the effectiveness of our optimizations, and compare the performance of the optimized method with the state-of-the-art on large-scale synthetic and real RDF datasets containing billions of triples. These experiments confirm that our SPAROL-UO query evaluation method is orders of magnitude more efficient than existing work.

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