An Extensible Framework for Query Optimization on TripleT-Based RDF Stores

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Resource description framework (RDF)

- W3C recommendation data model for linked data
- Graph-based data model, embodying lessons learned from previous data models
  - flexibly capture “things” and their relationships
  - don’t force artificial data/metadata distinctions
  - support full spectrum between structured and unstructured data
Resource description framework (RDF)

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  - flexibly capture “things” and their relationships
  - don’t force artificial data/metadata distinctions
  - support full spectrum between structured and unstructured data
- Uses web identifiers (URIs) to identify resources (“things”)
- Uses triples to state relationships between things
- serialization formats: N-triples, N3, RDF/XML
RDF datasets (or: RDF graphs) consist of a finite number of triples:

(Chell, worksAt, ApertureLabs)
(Chell, hasOccupation, "Test Subject")
(ApertureLabs, hasSlogan, "A Trusted Friend in Science")
(Caroline, worksAt, ApertureLabs)
(Caroline, supervises, Chell)

Each triple is of the shape \((s, p, o)\).
SPARQL is the W3C recommended language for querying RDF datasets.

PREFIX ex: <http://example.org/>
SELECT ?slogan
WHERE {
  ex:Chell ex:worksAt ?x.
  ex:Caroline ex:worksAt ?x.
  ?x ex:hasSlogan ?slogan.
}
We focus on the core of SPARQL: Basic Graph Patterns (BGPs).
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BGPs are constructed from Simple Access Patterns (SAPs):

\((?x, \text{worksAt}, ?y)\)

A conjunction of one or more SAPs makes up a BGP:

\{(?x, \text{worksAt}, ?y), (\text{?y, hasSlogan, ?z})\}
Results of BGP queries are sets of variable bindings:
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(?x, worksAt, ?y)

On our example dataset, we get:

{(?x : Chell),
  (?y : ApertureLabs)}
{(?x : Caroline),
  (?y : ApertureLabs)}
Results of BGP queries are sets of variable bindings:

\{ (Chell, worksAt, ?x),
    (?x, hasSlogan, ?y) \}

On our example dataset, we get:

\{ (?x : ApertureLabs),
    (?y : "A Trusted...") \}
Storing and querying RDF datasets is supported by RDF stores (also known as triple stores). As adoption of RDF grows, RDF stores have to deal with increasingly large datasets.
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Two basic storage alternatives

1. Flat file of triples, one per line
   - ... sorted on some subset of positions (e.g., by subject values)

2. Index data structure: maps a search key to a set of matching triples
   - B+tree, hash table
state of the art in the market: value-based indexing

- MAP (2007)
- HexTree (2008)
- TripleT (2009)
Storage: on the varieties of triple indexes

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still in the research lab: structure-based indexing

- group triples both by the values they share and by the *structure* they share in the graph (e.g., similar neighborhood topology)
- structure-based indexing also successful in XML data management
Let’s focus on value-based indexes

For graph $G$, let

\[
S(G) = \{ s | (s, p, o) \in G \}
\]
\[
P(G) = \{ p | (s, p, o) \in G \}
\]
\[
O(G) = \{ o | (s, p, o) \in G \}
\]
\[
\mathcal{A}(G) = S(G) \cup P(G) \cup O(G)
\]
**MAP**: index all permutations of S, P, and O

\[
SPO = \{s\#p\#o \mid (s, p, o) \in G\} \\
SOP = \{s\#o\#p \mid (s, p, o) \in G\} \\
PSO = \{p\#s\#o \mid (s, p, o) \in G\} \\
POS = \{p\#o\#s \mid (s, p, o) \in G\} \\
OSP = \{o\#s\#p \mid (s, p, o) \in G\} \\
OPS = \{o\#p\#s \mid (s, p, o) \in G\} 
\]
MAP: index all permutations of S, P, and O

... times six
**MAP**: index all permutations of S, P, and O

Most popular approach to triple storage in practice

- **RDF-3X**: open source triple store
- **Virtuoso**: open source and commercial industrial-strength triple store
- **Sesame/BigData**: open source and commercial industrial-strength triple store
HexTree: index all pairs of S, P, and O

\[
egin{align*}
SP & = \{ s\#p \mid (s, p, o) \in G \} \\
SO & = \{ s\#o \mid (s, p, o) \in G \} \\
PS & = \{ p\#s \mid (s, p, o) \in G \} \\
PO & = \{ p\#o \mid (s, p, o) \in G \} \\
OS & = \{ o\#s \mid (s, p, o) \in G \} \\
OP & = \{ o\#p \mid (s, p, o) \in G \} 
\end{align*}
\]
**HexTree**: index all pairs of S, P, and O

... times six
HexTree: partners share payload

\[ SP \rightarrow \{ o \in \mathcal{O}(G) \mid (s, p, o) \in G \} \]
\[ PS \rightarrow \{ o \in \mathcal{O}(G) \mid (s, p, o) \in G \} \]
MAP and HexTree

- **strengths**: support all native joins expressible in the SPARQL WHERE clause, using fast “merge” joins, with no constraints on “schema”
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- **weaknesses**
  - lots of (redundant) storage space
  - access to multiple data structures to perform joins
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- lots of (redundant) storage space
- access to multiple data structures to perform joins
- in general, weak data locality
TripleT (Fletcher and Beck, CIKM 2009)
- index $A(G)$
- just one data structure – a three-way triple tree
Storage: on the varieties of triple indexes

TripleT

Diagram showing a triangle with nodes labeled with triplets and arrows indicating connections.
Storage: on the varieties of triple indexes

TripleT

Diagram showing the components and structure of TripleT.
**TripleT** advantages

- retains strengths of MAP and HexTree
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Leads to significantly reduced storage and query processing costs
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Idea appears in various guises in products

- open source TU/e implementation (this work)
  
  https://github.com/b-w/TripleT

- open source 4store engine

- and storage of RDF in so-called column stores
  
  - open source MonetDB
  
  - commercial Vertica
Also, relational DB-based RDF solutions are available in industrial-strength IBM DB2 and Oracle Database 11G products.

Both use various hybrids of the MAP, HexTree, and TripleT indexes.
An RDF dataset consists entirely of triples. This results in queries with many self-joins, which provides a novel performance challenge.

Currently, BGP query optimization is an active area of research, where the challenge is to deliver interactive response times on BIG graphs.
Query optimization on TripleT has not been systematically studied up to this point.

This work: we focus on optimization of query plans for BGP queries on TripleT.

We investigate how the use of heuristics and external information (e.g. dataset statistics) can contribute towards a more intelligent way of generating query plans, minimizing query execution time over the TripleT RDF store.
Contributions

- We present an open source secondary-memory implementation of the TripleT RDF store and index.
- We propose a framework for query optimization in the form of a generic, rule-based algorithm used for generating query plans for given Basic Graph Pattern queries.
- We introduce a number of rules for use by our optimization framework.
- We implement this optimization framework on top of our TripleT store.
- We perform an extensive empirical study into the effectiveness of our rules in generating optimized query plans.
To answer a BGP query

\[(s_1, p_1, o_1) \land (s_2, p_2, o_2) \land \cdots \land (s_n, p_n, o_n)\]

we need to compute the natural join over the binding sets produced by its SAPs:

\[(s_1, p_1, o_1) \Join (s_2, p_2, o_2) \Join \cdots \Join (s_n, p_n, o_n)\]
Query optimization

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\[(s_1, p_1, o_1) \bowtie (s_2, p_2, o_2) \bowtie \cdots \bowtie (s_n, p_n, o_n)\]

Doing so requires a (physical) query plan.

---

Hash Join

Merge Join (?x)

P-Scan (?y, d, e)

S-Scan (a, b, ?x)

P-Scan (?x, c, ?y)
Deciding in what order to perform these joins is crucial to query response time.

Different plans for the same query can have response times several orders of magnitude apart.
Goal: separation of algorithm and rules.

Solution: a generic algorithm for query plan generation, containing several decision points which are guided by a set of rules.
The steps of the algorithm are:

1. Compute the *atom collapse* of a given BGP.
2. From the atom collapse, compute a *join graph*.
3. From the join graph, compute a query plan.
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Several rules work to obtain two primary goals:

1. Maximize merge joins
2. Minimize (expected) intermediate results
As an example, consider the following BGP:

$$(a, b, ?x)_1 \land (?x, c, ?y)_2 \land (?y, d, e)_3$$
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We then use this to obtain a join graph:
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\[(a, b, ?x)_1 \land (?x, c, ?y)_2 \land (?y, d, e)_3\]

The join graph is then used to obtain a query plan:
Seed rules

**Seed rule 1:** select one preferred seed for each distinct SAP in the input set based on the positions of the atoms in the SAP following the ordering \(s \succ o \succ p\).
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- **Seed rule 2**: prioritizes the atoms in the SAP according to their selectivity, as indicated by dataset statistics.
Join rules

- **Join rule 1**: merge join prioritization.
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- **Join rule 2**: prioritize joins involving the most selective SAPs
  
  \[(s, p, o) \succ (s, ?, o) \succ (s, p, ?) \succ (?, p, o) \succ (s, ?, ?) \succ 
  (?, ?, o) \succ (?, p, ?) \succ (?, ?, ?)\]

- **Join rule 3**: prioritize joins between SAPs that have the most selective positioning of join variables

- **Join rule 4**: prioritize joins between SAPs which feature a literal value (e.g. “Sue”) in one of its positions, over those featuring only URIs (e.g. “http://example.org/Sue”)

- **Join rule 5**: expected output minimization. Using a precomputed statistics database, output sizes of joins are known in advance.
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  selectivity of atoms, pairs of atoms, and joins of SAPs
Experiments: goals

Goal is to answer a number of questions:

1. What is the contribution of each of the rules?
2. How effective are combinations or subsets of rules?
3. What is the added value of statistics?
4. How do the rules perform under different datasets and queries?
9 different datasets, from 3 different sources:

- DBpedia
- SP²Bench
- UniProt

...and in 3 different sizes:

- 100,000 triples
- 1,000,000 triples
- 10,000,000 triples
Queries in different shapes:

...and of different sizes and selectivity.
We test with different configurations of rules.

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<th>Rules</th>
<th>S1</th>
<th>S2</th>
<th>J1</th>
<th>J2</th>
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Each of the runs has a particular purpose:

- **A-runs**: designed to test the heuristics rules against the statistics rules.
- **B-runs**: aim to get a sense of the value of the individual heuristics rules.
- **C-runs**: focus on the ordering of rules.
- **D-runs**: are used to test different subset combinations of rules.

For each configuration, we test on all datasets with all queries, and observe performance.
Experiments: setup

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For each configuration we test on all datasets with all queries, and observe performance.

The B- and C-runs confirmed the observations in the other runs, so we only show the main observations from the A- and D-runs
Experiments: overview of runs A-3, A-4, and D-2

**SP²Bench**

- **Dataset size (number of triples)**: 100K, 1M, 10M
- **Execution time (ms, logscale)**

**UniProt**

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- **Dataset size (number of triples)**: 100K, 1M, 10M
- *Excluding 15 / 60 data points with values > 1E+06
Experiments: A runs on SP2Bench

Run A-3

A BGP with 8 SAPs. Execution of A-4 is an order of magnitude slower.
Simple heuristics can consistently generate good query plans.

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The impact of statistics and the statistical prioritization join rule is measurable but limited.

Simple heuristic-driven rule sets such as A-3 and D-2 are the clear winners.
What’s next?

Future directions:

- Extensively study the problem of statistics-based query optimization on TripleT.
- Study runtime optimization strategies in our framework, such as sideways information passing.
- Continue open-source development of TripleT into a fully-featured RDF data management solution.
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Thank you!