

Greedy list intersection

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Abstract

A common technique for processing conjunctive queries is to first match each predicate separately using an index lookup, and then compute the intersection of the resulting row-id lists, via an AND-tree. The performance of this technique depends crucially on the order of lists in this tree: it is important to compute early the intersections that will produce small results. But this optimization is hard to do when the data or predicates have correlation.

We present a new algorithm for ordering the lists in an AND-tree by sampling the intermediate intersection sizes. We prove that our algorithm is near-optimal and validate its effectiveness experimentally on datasets with a variety of distributions.

1 Introduction

Correlation is a persistent problem for the query processors of database systems. Over the years, many have observed that the standard System-R assumption of independent attribute-value selections does not hold in practice, and have proposed various techniques towards addressing this (e.g, [8]).

Nevertheless, query optimization is still an unsolved problem when the data is correlated, for two reasons. First, the multidimensional histograms and other synopsis structures used to store correlation statistics have a combinatorial explosion with the number of columns, and so are very expensive to construct as well as maintain. Second, even if the correlation statistics were available, using correlation information in a correct way requires the optimizer to do an expensive numerical procedure that optimizes for maximum entropy [10]. As a result, most database implementations still rely heavily on independence assumptions.

1.1 Correlation problem in Semijoins

One area where correlation is particularly problematic is for semijoin operations that are used to answer conjunctive queries over large databases. In these operations, one separately computes the

set of objects matching each predicate, and then intersects these sets to find the objects matching the conjunction. We now consider some examples:

1. *Star joins*: The following query analyzes coffee sales in California by joining a fact table Orders with multiple dimension tables:

```
SELECT S.city, SUM(O.quantity), COUNT(E.name)
FROM orders O, cust C, store S, product P, employee E
WHERE O.cId=C.id and O.sId=S.id and O.pId=P.id and O.empld=e.id
      C.age=65, S.state=CA, P.type=COFFEE, E.type=TEMP
GROUP BY S.city
```

Many DBMSs would answer this query by first intersecting 4 lists of row ids (RIDs), each built using a corresponding index:

$$\begin{aligned} L_1 &= \{\text{Orders.id} \mid \text{Orders.cId} = \text{Cust.id}, \text{Cust.age} = 65\}, \\ L_2 &= \{\text{Orders.id} \mid \text{Orders.sId} = \text{Store.id}, \text{Store.state} = \text{CA}\}, \\ &\dots; \end{aligned}$$

and then fetching and aggregating the rows corresponding to the RIDs in $L_1 \cap L_2 \cap \dots$.

2. *Scans in Column Stores*: Recently there has been a spurt of interest in column stores (e.g, [15]). These would store a schema like the above as a denormalized “universal relation”, decomposed into separate columns for type, state, age, quantity, and so on. A column store does not store a RID with these decomposed columns; the columns are all sorted by RID, so the RID for a value is indicated by its position in the column. To answer the previous example query, a column store will use its columns to find the list of matching RIDs for each predicate, and then intersect the RID-lists.

3. *Keyword Search*: Consider a query for (“query” and (“optimisation” or “optimization”)) against a search engine. It is typically processed as follows. First, each keyword is separately looked up in an (inverted list) index to find 3 lists L_{query} , $L_{\text{optimisation}}$, and $L_{\text{optimization}}$ of matching document ids, and the second and third lists are merged into one sorted list. Next, the two remaining lists are intersected and the ids are used to fetch URLs and document summaries for display.

The intersection is often done via an *AND-tree*, a binary tree whose leaves are the input lists and whose internal nodes represent intersection operators. The performance of this intersection depends on the ordering of the lists within the tree. Intuitively, it is more efficient to form smaller intersections early in the tree, by intersecting together smaller lists or lists that have fewer elements in common.

Correlation is problematic for this intersection because the intersection sizes can no longer be estimated by multiplying together the selectivities of individual predicates.

1.2 State of the Art

The most common implementation of list intersection in data warehouses, column stores, and search engines, uses left-deep AND-trees where the k input lists L_1, L_2, \dots, L_k are arranged by increasing (estimated) size from bottom to top (in the tree). The intuition is that we want to form smaller intersections earlier in the tree. However, this method may perform poorly when the predicates are correlated, because a pair of large lists may have a smaller intersection than a pair of small lists.

Correlation is a well-known problem in databases and there is empirical evidence that correlation can result in cardinality estimates being wrong by many orders of magnitude, see e.g. [14, 8].

An alternative implementation proposed by Demaine et al [5] is a round-robin intersection that works on sorted lists. It starts with an element from one list, and looks for a match in the next list. If none is found, it continues in a round-robin fashion, with the next higher element from this second list. This is an extension to k lists of a comparison-based process that computes the intersection of two lists via an alternating sequence of doubling searches.

Neither of these two solutions is really satisfying. The first is obviously vulnerable to correlations. The second is guaranteed to be no worse than a factor of k from the best possible intersection (informally, because the algorithm operates in round-robin fashion, once in k tries it has to find a good list). But in many common inputs it actually performs a factor k worse than a naive left-deep AND-tree. For example, suppose the predicates were completely independent and selected rows with probabilities $p_1 \leq p_2 \leq \dots \leq p_k$, and suppose further that $\{p_j\}$ forms (or is dominated by) a geometric sequence bounded by say $1/2$. For a domain with N elements, an AND-tree that orders the lists by increasing size would take time $O(N(p_1 + p_1p_2 + \dots + p_1p_2 \dots p_{k-1})) = O(p_1N)$, while the round-robin intersection would take time proportional to $Nk/(\frac{1}{p_1} + \dots + \frac{1}{p_k}) = \Omega(kp_1N)$. This behavior was also experimentally observed in [6].

The round-robin method also has two practical limitations. First, it performs simultaneous random accesses to k lists. Second, these accesses are inherently serial and thus have to be low-latency operations. In contrast, a left-deep AND-tree accesses only two lists at a time, and a straightforward implementation of it requires random accesses to only one list. Even here, a considerable speedup is possible by dispatching a large batch of random accesses in parallel. This is especially useful when the lists are stored on a disk-array, or at remote data sources.

1.3 Contribution of this paper

We present a simple adaptive greedy algorithm for list intersections that solves the correlation problem. The essential idea is to order lists not by their marginal (single-predicate) selectivities, but rather by their conditional selectivities with respect to the portion of the intersection that has already been computed. Our method has strong theoretical guarantees on its worst case performance.

We also present a sampling procedure that computes these conditional selectivities at query run time, so that no enhancement needs to be made to the optimizer statistics.

We experimentally validate the efficacy of our algorithm and estimate the overhead of our sampling procedure by extensive experiments on a variety of data distributions.

To streamline the presentation, we focus throughout the paper on the data warehouse scenario, and only touch upon the extension to other scenarios in section 4.

1.4 Other Related Work

Tree-based RID-list intersection has been used in query processors for a long time. Among the earliest to use the greedy algorithm of ordering by list size was [11], who proposed the use of an AND-tree for accessing a single table using multiple indexes.

Round-robin intersection algorithms first arose in the context of AND queries in search engines. Demaine et al [5] introduced and analyzed a round-robin set-intersection algorithm that is based on a sequence of doubling searches. Subsequently, Barbay et al [2] have generalized the analysis

of this algorithm to a different cost-model. Heuristic improvements of this algorithm were studied experimentally on Google query logs in [6, 3]. A probabilistic version of this round-robin algorithm was used by Raman et al [13] for RID-list intersection.

In XML databases, RID-list intersection is used in finding all the matching occurrences for a twig pattern that selection predicates are on multiple elements related by an XML tree structure. [4] proposed a holistic twig join algorithm, *TwigStack*, for matching an XML twig pattern. IBM’s DB2 XML has implemented a similar algorithm for its XANDOR operator [1]. *TwigStack* is similar to round-robin intersection, navigating around the legs for results matching a pattern. Our algorithm can be applied to address correlation in all of these cases.

The analysis of our adaptive greedy algorithm uses techniques from the field of Approximation Algorithms. In particular, we exploit a connection to a different optimization problem, called the Min-Sum Set-Cover (MSSC) problem. In particular, we shall rely on previous work of Feige, Lovász and Tetali [7], who proved that the greedy algorithm achieves 4-approximation for this problem (see Appendix A.1 for details).

Pipelined filters. A variant of MSSC, studied by Munagala et al [12], is the *pipelined filters problem*. In this variant, a single list L_0 is given as the “stream” from which tuples are being generated. All predicates are evaluated by scanning this stream, so they can be treated as lists that support only a `contains()` interface that runs in $O(1)$ time. The job of the pipelined filters algorithm is to choose an ordering of these other lists. [12] apply MSSC by treating the complements of these lists as sets in a set covering. They show that the greedy set cover heuristic is a 4-approximation for this problem, and also study the online case (where L_0 is a stream of unknown tuples).

The crucial difference between this problem and the general list intersection problem is that an algorithm for pipelined filters is restricted to use a particular L_0 , and apply the other predicates via `contains()` only. Hence, every algorithm has to inspect every element in the universe at least once. In our context, this would be no better than doing a table scan on the entire fact table, and applying the predicates on each row. Another difference is in the access to the lists – our setting accommodates sampling and hence estimation of (certain) conditional selectivities, which is not possible in the online (streaming) scenario of [12], where it would correspond to sampling from future tuples. Finally, the pipeline of filters corresponds to a left-deep AND-tree, while our model allows arbitrary AND-trees; for example, one can form separate lists for say `age=65` and `type=COFFEE`, and intersect them, rather than applying each of these predicates one by one on a possibly much larger list.

2 Our Greedy Algorithm

Our list intersection algorithm builds on top of a basic infrastructure: the access method interface provided by the lists being intersected. The capability of this interface determines the cost model for intersection.

The two scenarios presented in the introduction – data warehouse and column stores, lead to different interfaces (and different cost models). In this section we present our intersection algorithm, focusing on the data warehouse scenario and associated cost model. We will return to the other scenario in Section 4.

2.1 List Interface

The lists being intersected are specified in terms of tables and predicates over the tables, such as $L_1 = \{\text{Orders.id} \mid \text{Orders.cId} = \text{Cust.id}, \text{Cust.age} = 65\}$. We access the elements in each such list via two kinds of operations:

- *iterate()*: This is an ability to iterate through the elements of the list in pipelined fashion.
- *contains(rid)*: This looks up into the list for occurrence of the specified RID.

These two basic routines can be implemented in a variety of ways. One could retrieve and materialize all matching RIDs in a hash table so that subsequent *contains()* is fast. Or, one could implement *contains()* via a direct lookup into the fact table to retrieve the matching row, and evaluate the predicate directly over that row.

2.2 The Min-Size Cost Model

Given this basic list interface, an intersection algorithm is implemented as a tree of pairwise intersection operators over the lists. The cost (running time) of this AND-tree is the sum of the costs of the pairwise intersection operators in the tree.

We model the cost of intersecting two lists $L_{\text{small}}, L_{\text{large}}$ (we assume wlog that $|L_{\text{small}}| \leq |L_{\text{large}}|$) as $\min\{|L_{\text{small}}|, |L_{\text{large}}|\}$. We have found this cost model to match a variety of implementation, that all follow the pattern of — *for each element x in L_{small} , check if L_{large} contains x* . Here, L_{small} has to support an iterator interface, while L_{large} need only support a *contains()* operation that runs in constant time.

We observe that under this cost model, the optimal AND-tree is always a left-deep tree (Claim A.1 in the appendix). Thus, L_{small} has to be formed explicitly only for the left-most leaf; it is available in a pipelined fashion at all higher tree levels.

In our data warehouse example, L_{small} is formed by index lookups, such as on *Cust.age* to find $\{\text{Cust.id} \mid \text{age} = 65\}$, and then for each *id* x therein, lookup an index on *Orders.cId* to find $\{\text{Orders.id} \mid \text{cId} = x\}$. L_{large} needs to support *contains()*, and can be implemented in two ways:

1. L_{large} can be formed explicitly as described for L_{small} , and then built into a hash table.
2. $L_{\text{large}}.\text{contains}()$ can be implemented without forming L_{large} , via index lookups: to check if a RID is contained in say $L_2 = \{\text{Orders.id} \mid \text{O.sId} = \text{Store.id} \text{ and } \text{Store.state} = \text{CA}\}$, we just fetch the corresponding row (by a lookup into the index on *Orders.id* and then into *Store.id*) and evaluate the predicate.

2.3 The Greedy Algorithm

We propose and analyze a new algorithm that orders the lists into a left-deep AND-tree, dynamically, in a greedy manner. Denoting the input k lists by $\mathcal{L} = \{L_1, L_2, \dots, L_k\}$, our greedy algorithm chooses lists G_0, G_1, \dots as follows:

1. Initialization: Start with a list of smallest size $G_0 = \operatorname{argmin}_{L \in \mathcal{L}} |L|$.
2. Iteratively for $i = 1, 2, \dots$: Assuming the intersection $G_0 \cap G_1 \cap \dots \cap G_{i-1}$ was already computed, choose the next list G_i to intersect with, such that the (estimated) size of $|(G_0 \cap G_1 \cap \dots \cap G_{i-1}) \cap (G_i)|$ is minimized.

We will initially assume that perfectly accurate intersection size estimates are available. In Section 3 we provide an estimation procedure that approximates the intersection size, and analyze the effect of this approximation on our performance guarantees.

One advantage of this greedy algorithm is its simplicity. It uses an left-deep AND-tree structure, similarly to what is currently implemented in most database systems. The AND-tree is determined only on-the-fly as the intersection proceeds. But this style of progressively building a plan fits well in current query processors, as demonstrated by systems like Progressive Optimization [9].

But a more important advantage of this greedy algorithm is that it attains worst-case performance guarantees, as we discuss next. For an input instance \mathcal{L} , let $\text{GREEDY}(\mathcal{L})$ denote the cost incurred by the above greedy algorithm on this instance, and let $\text{OPT}(\mathcal{L})$ be the minimum possible cost incurred by any AND-tree on this instance. We have the following result:

Theorem 2.1. *In the Min-Size cost model, the performance of the greedy algorithm is always within factor of 8 of the optimum, i.e., for every instance \mathcal{L} , $\text{GREEDY}(\mathcal{L}) \leq 8 \cdot \text{OPT}(\mathcal{L})$. Further, it is NP-hard to find an ordering of the lists that would give performance within factor better than $5/2$ of the optimum (even if the size of every intersection can be computed).*

The proof of Theorem 2.1 appears in Appendix A.2. We supplement our theoretical results with experimental evidence that our greedy algorithm indeed performs better than the commonly used heuristic, that of using a left-deep AND-tree with the lists ordered by increasing size. The experiments are described in detail in Section 5.

3 Sampling Procedure

We now revisit our assumption that perfectly accurate intersection size estimates are available to the greedy algorithm. We will use a simple procedure that estimates the intersection size within small absolute error, and provide rigorous analysis to show that it is sufficiently effective for the performance guarantees derived in the previous sections. As we will see in Theorem 3.4, the total cost (running time) of computing the intersection estimates is polynomial in k , the number of lists (which is expected to be small), and completely independent of the list sizes (which are typically large).

Proposition 3.1. *There is a randomized procedure that gets as input $0 < \varepsilon, \delta < 1$ and two lists, namely a list A that supports access to a random element, and a list B that supports the operation `contains()` in constant time; and produces in time $O(\frac{1}{\varepsilon^2} \log \frac{1}{\delta})$ an estimate s such that*

$$\Pr \left[s = |A \cap B| \pm \varepsilon |A| \right] \geq 1 - \delta.$$

Proof. (Sketch) The estimation procedure works as follows:

1. Choose independently $t = \frac{1}{\varepsilon^2} \log \frac{1}{\delta}$ elements from A ,
2. Compute the number s' of these elements that belong to B .
3. Report the estimate $s = \frac{s'}{t} \cdot |A|$.

The proof of the statement is a straightforward application of Chernoff bounds. □

In practice, the sampling step 1 can be done either by materializing the list A and choosing elements from it, or by scanning a pre-computed sample of the fact table and choosing elements that belong to A .

We next show that in our setting, the absolute error of Proposition 3.1 actually translates to a relative error. This relative error does not pertain to the intersection, but rather to its complement, as seen in the statement of the following proposition. Indeed, this is the form which is required to extend the analysis of Theorem 2.1.

Proposition 3.2. *Let $\mathcal{L} = \{L_1, \dots, L_k\}$ be an instance of the list intersection problem, and denote $I = L_1 \cap L_2 \cdots \cap L_j$. If $|I \cap L_m|$ is estimated using Proposition 3.1 for each $m \in \{j+1, \dots, k\}$ and m^* is the index yielding the smallest such estimate, then $|I \setminus L_{m^*}| \geq (1 - 2k\varepsilon) \max_m |I \setminus L_m|$.*

Proof. For $m \in \{j+1, \dots, k\}$, the estimate for $|I \cap L_m|$ naturally implies an estimate for $|I \setminus L_m|$, which we shall denote by s_m . By our accuracy guarantee, for all such m

$$s_m = |I \setminus L_m| \pm \varepsilon |I|. \quad (1)$$

Let m_0 be the index that really maximizes $|I \setminus L_m|$. Let m^* be the index yielding the smallest estimate for $|I \cap L_m|$, i.e., the largest estimate for $|I \setminus L_m|$. Thus, $s_{m^*} \geq s_{m_0}$, and using the accuracy guarantee (1) we deduce that

$$|I \setminus L_{m^*}| \geq s_{m^*} - \varepsilon |I| \geq s_{m_0} - \varepsilon |I| \geq |I \setminus L_{m_0}| - 2\varepsilon |I|. \quad (2)$$

The following lemma will be key to completing the proof of the proposition.

Lemma 3.3. *There exists $m \in \{j+1, \dots, k\}$ such that $|I \setminus L_m| \geq |I|/k$.*

Proof of Lemma 3.3. Recall that $\cap_i L_i = \emptyset$. Thus, every element in I does not belong to at least one list among L_{j+1}, \dots, L_k , i.e., $I \subseteq \cup_{m=j+1}^k (I \setminus L_m)$. By averaging, at least one of these lists $I \setminus L_m$ must have size at least $|I|/(k-j)$. \square

We now continue proving the proposition. Using Lemma 3.3 we know that $|I \setminus L_{m_0}| \geq |I|/k$, and together we conclude, as required, that

$$|I \setminus L_{m^*}| \geq |I \setminus L_{m_0}| - 2\varepsilon |I| \geq |I \setminus L_{m_0}| \cdot (1 - 2k\varepsilon).$$

This completes the proof of Proposition 3.2. \square

What accuracy (value of $\varepsilon > 0$) do we need to choose when applying Proposition 3.2 to the greedy list intersection? A careful inspection of the proof of Theorem 2.1 reveals that our the performance guarantees continue to hold, with slightly worse constants, if the greedy algorithm chooses the next list to intersect with using a constant approximation to the intersection sizes. Specifically, suppose that for some $0 < \alpha < 1$, the greedy algorithm chooses lists G_0, G_1, \dots (in this order), and that at each step j , the list G_j is only factor α approximately optimal in the following sense (notice that the factor α pertains not to the intersection, but rather to the complement):

$$|(G_0 \cap \dots \cap G_{j-1}) \setminus G_j| \geq \alpha \cdot \max_{L \in \mathcal{L}} |(G_0 \cap \dots \cap G_{j-1}) \setminus L|. \quad (3)$$

Note that $\alpha = 1$ corresponds to an exact estimation of the intersections, and the proof of Theorem 2.1 holds. For general $\alpha > 0$, an inspection of the proof shows that the performance guarantee of Theorem 2.1 increases by a factor of at most $1/\alpha$.

From Proposition 3.2, we see that choosing the parameter ε (of our estimation procedure) to be of the order of $1/k$ gives a constant factor approximation of the above form. Choosing the other parameter δ carefully gives the following:

Theorem 3.4. *Let every intersection estimate used in the greedy algorithm on input \mathcal{L} be computed using Proposition 3.1 with $\varepsilon \leq 1/(8k)$ and $\delta \leq 1/k^2$.*

(a) The total cost (running time) of computing intersection estimates is at most $O(k^4 \log k)$, independently of the list sizes.

(b) With high probability, the bounds in Theorem 2.1 hold with larger constants.

Proof. Part (a) is immediate from Proposition 3.1, Since the greedy algorithm performs at most k iterations, each requiring at most k intersection estimates. It thus remains to prove (b).

For simplicity, we first assume that the input instance $\mathcal{L} = \{L_1, \dots, L_k\}$ satisfies $\cap_i L_i = \emptyset$. By Propositions 3.2 and our choice of ε and δ , we get that, with high probability, every list G_j chosen by greedy at step j is factor $1/2$ approximately optimal in the sense of (3). It is enough that each estimate has, say, accuracy parameter $\varepsilon = 1/(8k)$ and confidence parameter $\delta = 1/k^2$.

Now consider a general input \mathcal{L} and denote $I^* = \cap_i L_i$. We partition the iterations into two groups, and deal with each group separately. Let $j' \geq 1$ be the smallest value such that $|G_0 \cap G_1 \cap \dots \cap G_{j'}| \leq 2|I^*|$. For iterations $j = 1, \dots, j' - 1$ (if any) we can essentially apply an argument similar to Proposition 3.2: we just ignore the elements in I^* , which are less than half the elements in $I = G_0 \cap G_1 \cap \dots \cap G_j$, and hence the term $1 - 2k\varepsilon$ should only be replaced by $1 - 4k\varepsilon$; as argued above, it can be shown that the algorithm chooses a list that is $O(1)$ -approximately optimal.

For iterations $j = j', \dots, k - 1$ (if any), we compare the cost of our algorithm with that of a greedy algorithm that would have had perfectly accurate estimates: Our algorithm has cost at most $O(|I^*|)$ per iteration, regardless of the accuracy of its intersection estimates, while if the estimates were perfectly accurate, would still cost at least $|I^*|$; hence, the possibly inaccurate estimates can increase our upper bounds by at most a constant factor. \square

4 Other Scenarios and Cost Models

We now turn to alternative cost model proposed by [5], which assumes that all the lists to be intersected have already been sorted. The column store example of Section 1.1 fits well into this model, because every column is kept sorted by RID.

On the other hand, this model is not valid in the data warehouse scenario because the lists are formed by separate index lookups for each matching dimension key. E.g., the list of RIDs matching `Cust.age = 65` is formed by separate lookups into the index on `Orders.cId` for each `Cust.id | age = 65`. The result of each lookup is individually sorted on RID, but the overall list is not.

4.1 The Comparisons Cost Model

In this model, the cost of intersecting two lists L_1, L_2 is the minimum number of comparisons needed to “certify” the intersection. This model assumes that both lists are already sorted by RID. Then, the intersection is computed by an alternating sequence of doubling searches¹ (see Figure 1 for illustration):

1. Start at the beginning of L_1 .
2. Take the next element in L_1 , and do a doubling search for a match in L_2 .
3. Take the immediately next (higher) element of L_2 , and search for a match in L_1 .
4. Go to Step 2.

¹ By doubling search we mean looking at values that are powers of two away from where the last search terminated, and doing a final binary search. We approximate this cost as $O(1)$.

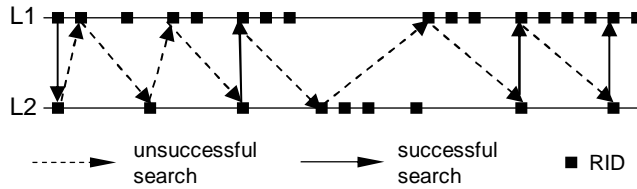


Figure 1: Intersecting two sorted lists

The number of searches made by this algorithm could sometimes be as small as $|L_1 \cap L_2|$, and at other times as large as $2 \min\{|L_1|, |L_2|\}$, depending on the “structure” of the lists (again approximating the cost of a doubling search as a constant).

4.2 The Round-Robin Algorithm

Demaine et al [5] and Barbay and Kenyon [2] have analyzed an algorithm that is similar to the above, but runs in a round-robin fashion over the k input lists. Their cost model counts comparisons, and they show that the worst-case running time of this algorithm is always within a factor of $O(k)$ of the smallest number of comparisons needed to certify the intersection. They also show that a factor of $\Omega(k)$ is necessary: there exists a family of inputs, for which no deterministic or randomized algorithm can compute the intersection in less than k times the number of comparisons in the intersection certificate.

4.3 Analysis of the Greedy algorithm

For the Comparisons model, we show next that the greedy algorithm is within a constant factor of the optimum plus the size of the smallest list, $\ell_{min} = |G_0| = \min_{L \in \mathcal{L}} |L|$; namely, for every instance \mathcal{L} , $\text{GREEDY}(\mathcal{L}) \leq 8 \cdot \text{OPT}(\mathcal{L}) + 16\ell_{min}$. We get around the factor $\Omega(k)$ lower bound of Barbay and Kenyon [2] by restricting OPT to be an AND-tree, and by allowing an additive cost based on ℓ_{min} (but independent of k). We further discuss the merits of the above bound vs. the factor $O(k)$ bound for round-robin in Section 6. We also prove that the above bound is the best possible, up to constant factors, since there are instances \mathcal{L} for which $\text{OPT}(\mathcal{L}) = O(1)$ and $\text{GREEDY}(\mathcal{L}) \geq (1 - o(1))\ell_{min}$. This instance shows that with our limited lookahead (information about potential intersections), paying $\Omega(\ell_{min})$ is essentially unavoidable, regardless of $\text{OPT}(\mathcal{L})$.

Theorem 4.1. *In the comparison cost model, the performance of the greedy algorithm is always within factor of 8 of the optimum (with an additive factor), i.e., for every instance \mathcal{L} , $\text{GREEDY}(\mathcal{L}) \leq 8 \text{OPT}(\mathcal{L}) + 16\ell_{min}$, where ℓ_{min} is length of the smallest input list. Further, there is a family of instances \mathcal{L} for which $\text{GREEDY}(\mathcal{L}) \geq (1 - o(1))(\text{OPT}(\mathcal{L}) + \ell_{min})$.*

The proof is in Appendix B. Note that the optimum AND-tree for the Min-Size model need not be optimal for the Comparison model and vice versa. Also note that if we only have estimates of intersection sizes (using Proposition 3.1), the theoretical bounds hold, with slightly worse constants.

5 Experimental Results

We validate the practical value of our algorithm via an empirical evaluation that addresses two questions:

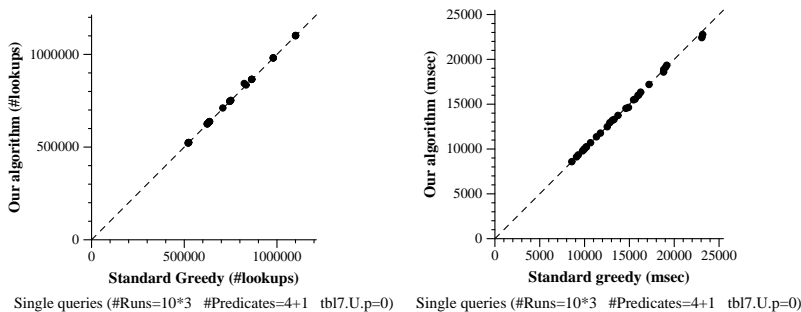


Figure 2: No degradation in performance, even with no correlations.

1. How does the algorithm perform in the presence of correlations? In particular, is it really more resilient to correlations than the common heuristic, and if so, is the performance improvement substantial?
2. Is the sampling overhead negligible? Specifically, under what circumstances can this algorithm be less efficient than the common heuristic, and by how much.

To answer these question we conducted experiments that compare the performance of the two algorithms in various synthetic situations, including a variety of data distributions (especially different correlations and soft functional dependencies), and various query patterns. The theoretical analysis suggests that our algorithm will outperform the common heuristic by picking up correlations in the data, namely, by avoiding positive correlations and embracing negative correlations. Furthermore, the sampling procedure has been shown to use very few samples, and is thus expected to have a rather negligible impact when the data is even moderately large. However, our theoretical analysis only counts certain operations and thus ignores a multitude of important implementation issues, like disk and memory operations, together with their access pattern (sequential/random) and their caching policies, and so forth.

5.1 Basic Setup

We implemented (a simple variant of) the first example mentioned in the Introduction, namely, conjunctive queries over a fact table. In all the experiments, there is a fact table with 10^7 rows, each identified by a unique row id (RID). The fact table has 8 columns, denoted by the letter A to H , that are partitioned into 3 groups: (1) $A - E$, (2) $F - G$, and (3) H ; columns from different groups are generated independently, and columns within a group are correlated in a controlled manner, as they are generated by soft functional dependencies. The fact table is accompanied by an index for each of the columns. All our queries are conjunctions of 5 predicates, where each predicate is specified by an attribute and a range of values, e.g. $B \in [7, 407]$. Each predicate in our experiments typically has selectivity of 10%, but their ranges may be correlated. It is important to note that both the fact table and the indices were written onto disk, and all query procedures rely on direct access to the disk (but with a warm start). A more detailed description of the data and queries generation is given below.

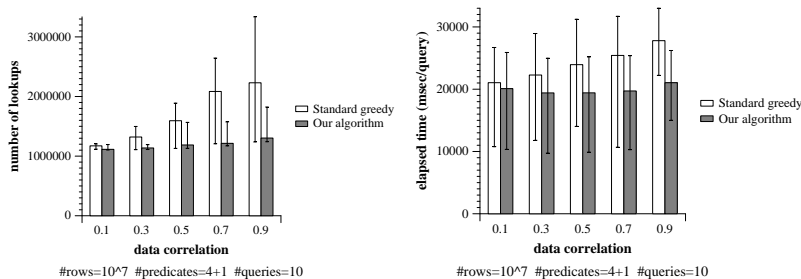


Figure 3: Our algorithm offers improved performance by detecting and exploiting correlations.

5.2 Machine Info

All experiments were carried out on an IBM POWER4 machine running AIX 5.2. The machine has one CPU and 8Gb of internal memory and, unless stated otherwise, the data (fact tables and indexes) was written on a local disk. We note however that our programs perform a query using relatively small amount of memory, and rely mostly on disk access.

5.3 Data Generation

As mentioned above, our fact table has 10^7 rows and 8 columns. Its basic setup is as follows (some variations will be discussed later): For each row, attribute A was generated uniformly at random from the domain $\{1, 2, \dots, 10^5\}$. Then, attributes B through E were each generated from the previous one using a soft functional dependency: with probability $p \in [0, 1]$ its value was chosen to be a deterministic function of the previous attribute, and with probability $1 - p$ its value was uniform at random. Attributes B, C, D have domain sizes $10^4, 10^3$, and 10^2 , respectively, and for each one the functional dependency is simply the previous attribute divided by 10 (i.e., $B = A/10$, $C = B/10$ and $D = C/10$). Attribute E has the same domain as attribute A and its functional dependency is that of complementarity, i.e. $E = 10^5 - A$. Attributes $F - G$ were generated similarly to $A - B$, but independent of A and B , and attribute H was generated similarly to A but again independently of all others. This model of correlations aims to capture common instances; for example, a car’s make (e.g. Toyota) is often a function of the car’s model (in this case, Camry, Corolla, etc.). The single parameter p gives us a convenient handle on the amount of correlation in the data.

Our queries are all a conjunction of 5 predicates, for example

$$\begin{aligned}
 A \in [1500, 2499] \quad \wedge \quad B \in [150, 200] \quad \wedge \quad C \in [15, 20] \\
 \wedge \quad E \in [2001, 3000] \quad \wedge \quad F \in [1, 999]
 \end{aligned}$$

which leads to various selectivities and correlations, depending on the choice of attributes. For example, in this query the first predicate have selectivity 1%, and the third one has selectivity 0.6%. Furthermore, attributes A and B are positively correlated, attributes A and E are negatively correlated, and attributed A and F are not correlated. Note that the correlations take effect in the query only if their ranges are compatible. It’s important to note that incompatible ranges are less likely to occur, as it corresponds (continuing our earlier example) to querying abouts cars whose model is Civic and their make is Toyota (rather than Honda).

We need a methodology for repeating the same (actually similar) query several times, so that our experimental evaluation can report the average/maximum/minimum runtime for processing a query. We do this by using the same query “structure” but with different values. For instance, in the query displayed above, the range of A could be chosen at random, but then the ranges of B and C are derived from it as a linear function. Whenever we report an aggregate for $n > 1$ repetitions of a query, we use the notion of a *warm start*, which means that $n + 1$ queries are run, one immediately after the other, and the aggregate is computed on all but the first query.

5.4 Experiments on Independent Data

The first experiment is designed to tackle the second question raised at the beginning of Section 5. We compare the performance of our algorithm with that of the common heuristic in the absence of any correlations in data. In this experiment, the attributes are generated independently of each other, by setting the correlation parameter p to 0. Our algorithm will expend some amount of computation and data lookups towards estimating intersection sizes (via sampling), but we expect it will find no correlations to exploit. The results are depicted in Fig. 2. The chart on the left measures the number of *table lookups*, i.e. calls to the method `contains()`, and the one of the right measures the total elapsed time measured in milliseconds. (The former is the main ingredient of our theoretical guarantees.) Each point in the plot represents a run of the same query under the two algorithms; the point’s x coordinate represents the performance of the common heuristic, and its y coordinate represents the performance of our algorithm. We generated 30 points by running 10 different queries with predicate selectivities ranging from 1/10 to 1/20 (for all predicates), and repeating each such query 3 times at random. The predicates used were $A - D$ and F (but it should not matter in this plot because $p = 0$).

Notice that all points are very close to the line $y = x$ (i.e. at 45 degrees), shown by a dashed line. This is true both in terms of lookups operations and in terms of elapsed time. Thus, the two algorithms have essentially the same performance, showing that the sampling overhead is negligible.

5.5 Experiments on Correlated Data

The next experiment is designed to compare our algorithm with the common heuristic in the presence of correlation in data. The data is generated according to the distribution described in Section 5.3, with the correlation parameter p varying from 10% to 90%. The results are depicted in Fig 3. The chart on the left shows the number of *table lookups*, i.e. calls to the method `contains()`, and the one of the right shows the total elapsed time measured in milliseconds. (The former is the main ingredient of our theoretical guarantees.) The wide bars represent the average over 10 queries, and the thin lines plot the entire range (minimum to maximum). All queries had the same structure, namely 4 correlated attributes ($A - D$) and 1 independent attribute (E). The selectivity of each of these 5 predicates is the same, 10%.

Observe that our algorithm consistently performs better than the common heuristic even in the presence of moderate correlations. The improvement offered by our algorithm becomes more substantial as the correlation p increases. In fact, the performance of the common heuristic degrades as the correlation increases, while our algorithm maintains a rather steady performance level— this is strong evidence that our algorithm successfully detects and overcomes the correlations in the data. This gives a positive answer to the first question raised at the beginning of Section 5, at least

in a basic setup; later, we will examine this phenomenon under other variants of correlations in the data and queries.

The improvement in elapsed time is smaller than in the number of lookups; for example, at $p = 0.9$, the average number of lookups per query decreases by 42%, and the average elapsed time decreases by 25%. The reason is that the runtime is affected by the overhead of iterating over the lists (especially the first one); this overhead tends to be a small, but non-negligible, portion of the computational effort.

Finally, the errors in both algorithms are mostly similar, except that the common heuristic tends to have minima that are smaller (when compared to the respective average). The reason is simple—the heuristic essentially chooses a random ordering of the predicates (since they all have the same selectivity) and thus it every once in a while it happens to be lucky in choosing an ordering that exploits the correlations.

Varied Selectivity. In this experiment we test the dependence of our results above on the selectivity of the predicates (i.e. on the size of the lists). Recall that in the base experiment we used a standard selectivity parameter to get lists which are about 10% of the domain size, for each attribute A, B, C, D and F . Here we increase the selectivity of predicates B, C and D in the following manner: the selectivity of B, C and D go up to 0.8^{-1} , 0.8^{-2} and 0.8^{-3} respectively, causing the corresponding list sizes to shrink by factors of 0.8 , 0.8^2 and 0.8^3 respectively. We observe (see Fig 4) that our algorithm continues to give improved performance over greedy, and essentially with the same factors as before (the only difference being that now both algorithms perform faster than before, since the list sizes are smaller).

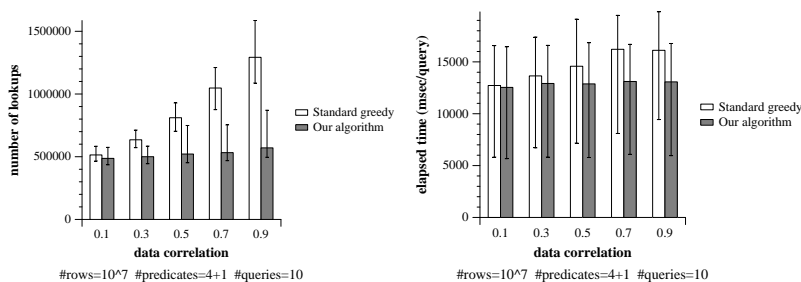


Figure 4: A comparison using predicates of higher selectivity.

Negative Correlation. In the next experiment we confirm the intuition that our algorithm does well in leveraging negative correlations in data. Here we replace the predicate based on attribute F by a predicate based on attribute E . Recall that F was independent of A , while E is negatively correlated with A , via the soft functional dependency of inversion. We observe in the experiments that the algorithm continues to perform much better than the standard greedy, both in terms of running time and number of lookups (see Fig 5). In fact we also observe that with a high correlation parameter, the improvement in performance obtained by our algorithm is much greater than in the case when we had only positive correlations. Indeed, the algorithm seems to find the correct negatively correlated predicates (A and E) to intersect in the first step, which immediately reduces the size of the intersection by a large amount, and results in savings later too. We see that in the case of very high correlations (90%), the standard greedy algorithm takes about 1.35

times as much processing time on the average, and makes about 2.5 times as many lookups as our algorithm on the average. Also we observe that the spread in the processing time and in the number of lookups is consistently smaller for our algorithm, with the spread being close to 0 as the correlation parameter increases. With correlation 90%, the maximum number of lookups made by the standard greedy is almost 3.5 times those made by our algorithm.

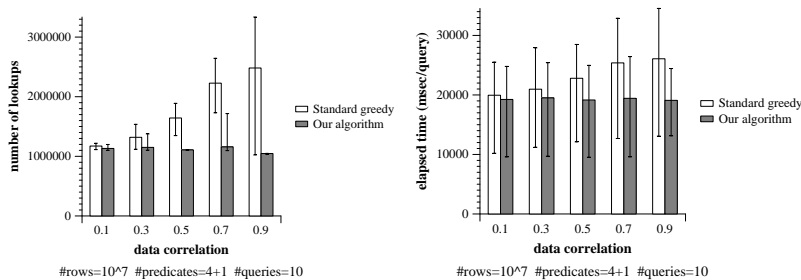


Figure 5: Leveraging negative correlations in data.

Skewed Data. In the next set of experiments we change the distribution from which the data is drawn. Instead of a uniform distribution, as in the previous experiments, now we pick our data from a very skewed distribution, namely a Zipf (Power-Law) Distribution. Again, we vary our correlation parameter from 10% to 90%. Our experimental results (see Fig 6) show that the performance of our algorithm remains significantly better than that of the common heuristic, confirming the hypothesis that the improved performance comes from correlations in data and is independent of the base distribution.

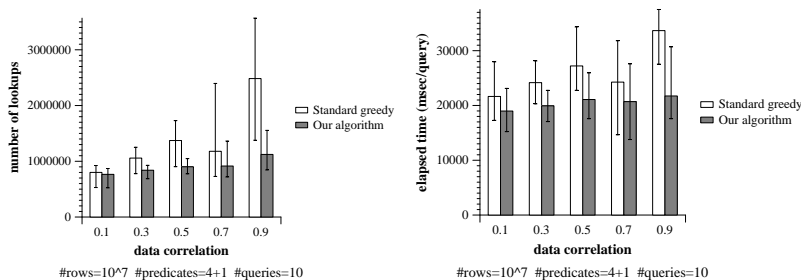


Figure 6: Data generated from a Zipf Distribution.

Sample Size. Next, we test the dependence of our algorithm on the number of samples it takes in order to estimate the intersection sizes. Theoretically we have guaranteed that a small sample size (polynomial in k , the number of lists) suffices to get good enough estimates. In the previous experiments we have verified that indeed such small samples can lead to significant improvement in efficiency, via our algorithm. In this experiment, we observe the dependence on the sample size, by varying the sample size from 1/128 to 64 times the standard value. As expected, Fig. 7 shows that when the sample size is too small, the estimates are not good enough, and the maximum as well as the average times and number of lookups are large. In fact, the smallest run times and lookups occur when the sample size factor is the one used in our previous experiments.

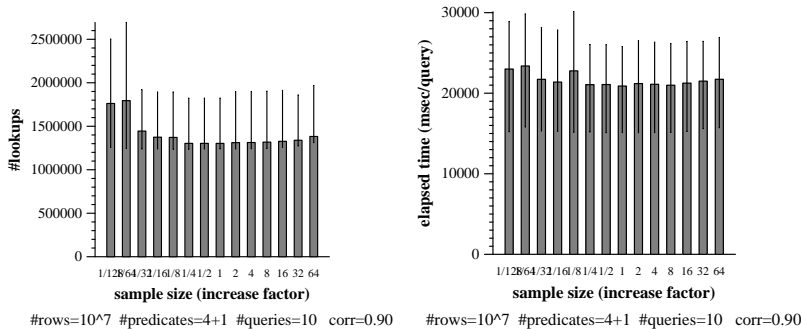


Figure 7: The dependence on the sample size used for intersection estimation

Latency. Finally, we evaluate the impact of latency (in access to the data) on the performance of the two algorithms. As pointed out earlier, larger latency may arise for various reasons such as storage specs (e.g. disk arrays) or when combining data sources (e.g. over the web). We thus compare the performance of the algorithms in the usual configuration when the data resides on a local disk, with one where the data resides on a network file system. Although the number of lookups is the same in both experiments, Fig. 8 shows that our improvement in the elapsed time becomes bigger when data has high latency (resides over the network). The reason is that high latency has more dramatic effect on random access to the data (lookup operations) vs. sequential access (iterating over a list).

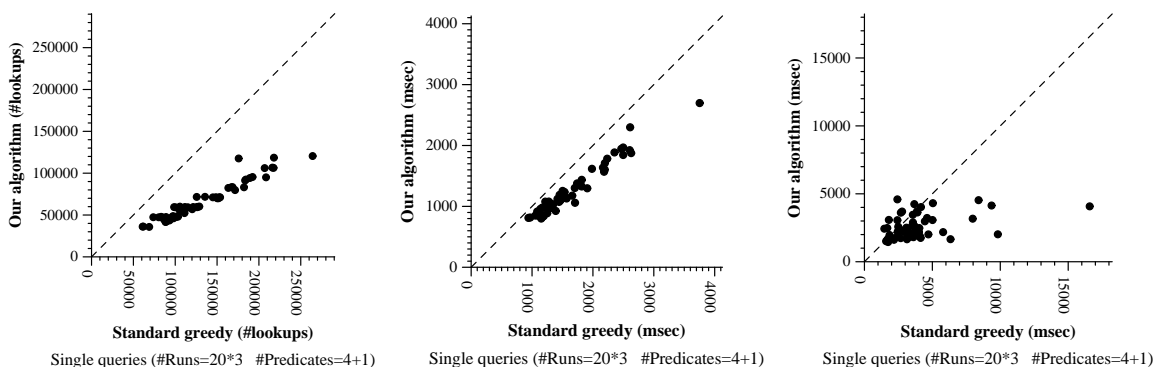


Figure 8: Comparison between elapsed time using a local disk (middle chart) and over the network (right). The number of lookups (left chart) is independent of implementation.

6 Conclusions and Future Work

We have proposed a simple greedy algorithm for the list intersection problem. Our algorithm is similar in spirit to the most commonly used heuristic, which orders lists in a left-deep AND-tree in order of increasing (estimated) size. But in contrast, our algorithm does have provably good worst-case performance, and much greater resilience to correlations between lists. In fact, the intuitive qualities of that common heuristic can be explained (and even quantified analytically) via our analysis: if the lists are not correlated, then list size is a good proxy for intersection size, and hence our analysis still holds. Our experimentation shows that these features of the algorithm

indeed occur empirically, under a variety of of data distributions and correlations. In particular, the overhead of our algorithm is in requiring new estimates after every intersection operation; overall, this is factor k more overhead than the heuristic approach, but as we showed, this cost is worth paying in most practical scenarios.

It is natural to ask for an algorithm whose running time achieves the “best of both worlds” (without running both the round-robin and the greedy algorithms in parallel until one of them terminates). Demaine et al [6] evaluate a few hybrid methods, but it seems that this question is still open. Another important question is whether these techniques can be used to do join ordering.

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A Proofs: Min Size Cost Model

A.1 The Min-Sum Set-Cover Problem

The approximation guarantees of our adaptive greedy algorithm exploit a connection to a different optimization problem, called the *min-sum set-cover* problem, and abbreviated in the sequel to MSSC. In this problem, the input is a set of lists $\mathcal{L} = \{L_1, L_2, \dots, L_m\}$, and the goal is to pick a sequence of lists $L_{i_1}, L_{i_2}, \dots, L_{i_t}$ from \mathcal{L} that covers all elements in the universe $U = \cup_{i=1}^m L_i$ so as to minimize the objective function $\sum_{e \in U} \min\{j | e \in L_{i_j}\}$ (in other words, we measure the time required to cover every element and sum over all elements).

Consider the following greedy algorithm for MSSC, which we denote by $\text{GREEDY}_{\text{MSSC}}$: Initially pick the list with the largest number of elements from \mathcal{L} , and from there on pick the list that covers the maximum number of yet uncovered elements. Feige, Lovász and Tetali [7] showed the following (worst-case) results.

Theorem A.1 ([7]). $\text{GREEDY}_{\text{MSSC}}$ achieves 4 approximation for the MSSC problem.

Theorem A.2 ([7]). For every $\varepsilon > 0$, it is NP-Hard to approximate the MSSC problem to within a factor of $4 - \varepsilon$.

For $0 < \alpha \leq 1$, let $\text{GREEDY}_{\text{MSSC}}^\alpha$ be a variant of the greedy algorithm that chooses at every step a list that is within factor α of optimal. Specifically, the list $L_j^\alpha \in \mathcal{L}$ that is picked at step $j = 1, 2, \dots$, satisfies the following condition:

$$|(U \setminus (L_1^\alpha \cup \dots \cup L_{j-1}^\alpha)) \cap L_j^\alpha| \geq \alpha \cdot \max_{L \in \mathcal{L}} |(U \setminus (L_1^\alpha \cup \dots \cup L_{j-1}^\alpha)) \cap L|.$$

Note that $\text{GREEDY}_{\text{MSSC}}$ is exactly the algorithm $\text{GREEDY}_{\text{MSSC}}^1$. By inspection of [7], we see that the proof of Theorem A.1 can be generalized as follows.

Theorem A.3. Let $0 < \alpha \leq 1$. Then $\text{GREEDY}_{\text{MSSC}}^\alpha$ achieves $4/\alpha$ approximation for the MSSC problem.

A.2 Proof of Theorem 2.1

We proceed to prove Theorem 2.1, namely, analyze the performance of the greedy algorithm in the Min-Size cost model. Let the cost of intersecting two lists A and B be denoted by $c_{\text{minsize}}(A, B) = \min\{|A|, |B|\}$.

Proof of Theorem 2.1. Let the input \mathcal{L} consist of the lists L_1, L_2, \dots, L_k . We may assume that the optimal AND-tree for \mathcal{L} is left-deep, by using the following claim.

Claim A.1. There is an optimal AND-tree for \mathcal{L} (i.e., tree of minimum cost) that happens to be a left-deep tree.

Proof of Claim A.1. Given an AND-tree, we will use left spine to denote the set of internal vertices that can be reached from the root by traversing successive “left child” edges. Given a tree T , we will use $\text{LS}(T)$ to denote the number of internal nodes *not* in the left spine. By a simple rearrangement, one can always assume that for every node in the left spine, the left child of the node corresponds to

the smaller of the two lists that are intersected at the node. For the rest of the proof, we will assume this is true for all AND-trees that we consider. We will prove the result by induction on $\text{LS}(T)$. Let k denote the number of input lists (or leaves in the tree). Note that if $\text{LS}(T) = 0$ then the tree is already left-deep. Assume that the result holds for all optimal AND-trees T' such that $\text{LS}(T') \leq s$, where $s \leq k - 1$. Now let T be an optimal AND-tree such that $\text{LS}(T) = s + 1$. Now as $\text{LS}(T) \geq 1$, there exist three (intermediate) lists A , B and C (note that each of them might correspond to a sub-tree of T) such that the optimal algorithm corresponding to T intersects B and C first (denote the node by v_1) and then intersects their result with A (denote the node by v_2). Further, v_2 is in the left spine of T and v_1 is the right child of v_2 (and is thus, not in the left spine). Note that the cost paid at node v_2 is $\min(|A|, |B \cap C|)$, while the cost paid at node v_1 is $\min(|B|, |C|)$. By the assumption that the left child of a node in the left spine corresponds to the smaller list, we have $|A| \leq |B \cap C|$. Now while keeping the rest of T the same, rearrange the tree such that first A and B are intersected (call this node v'_2) and then their intersection is intersected with C (call this node v'_1) such that v'_2 is the left child of v'_1 , which in turn is the left child of the parent of the (original node) v_2 . Call this new tree T'' . Now the cost paid at node v'_2 is $\min(|A|, |B|)$, which is the same as the cost paid originally at node v_2 (note that $|A| \leq |B \cap C| \leq |B|$). Further, the cost paid at node v'_1 is $\min(|A \cap B|, |C|)$, which is at most the price paid originally at node v_1 (note that $|A \cap B| \leq |B|$). Thus, the overall cost of T'' is at most the cost of T . Further, $\text{LS}(T'') = \text{LS}(T) - 1 = s$ and it is easy to check that all internal nodes in the left spine of T'' have the smaller list as their left child. This proves the inductive step and hence, the proof is complete. \square

We will first look at the case when the intersection of all the lists is empty. W.l.o.g. assume that the optimal algorithm intersects lists from \mathcal{L} according to the sequence (L_1, L_2, \dots, L_t) . Further, let the greedy sequence be (G_0, \dots, G_s) , where for every $0 \leq i \leq s$, $G_i \in \mathcal{L}$. Note that $\bigcap_{i=1}^t L_i = \bigcap_{i=0}^s G_i = \emptyset$. Let OPT be the optimal cost, and let GREEDY be the cost of the greedy algorithm. Consider also the following modification of the optimal sequence: $(G_0, L_1, L_2, \dots, L_t)$ and call its cost OPT' .

Claim A.2. $\text{OPT}' \leq 2 \text{OPT}$

Proof. OPT' is a sequence of t intersections and OPT is a sequence of $t - 1$ intersections. The first intersection in OPT' has cost $|G_0|$ (since $|G_0| \leq |L_1|$). For $i = 2, \dots, t$, the i th intersection in OPT' has cost at most the cost of the $(i - 1)$ th intersection in OPT . This is because $c_{\text{minsize}}(G_0 \cap L_1 \cap L_2 \dots \cap L_{i-1}, L_i) \leq c_{\text{minsize}}(L_1 \cap L_2 \dots \cap L_{i-1}, L_i)$. Thus we have $\text{OPT}' \leq |G_0| + \text{OPT}$. Now, even the first intersection in OPT has cost $\min\{|L_1|, |L_2|\} \geq |G_0|$. The claim follows. \square

Consider now the following instance \mathcal{I} of the MSSC problem. The universe is G_0 , and the sets are $S_i = G_0 \setminus L_i$, $i = 1, \dots, k$. Note that $\bigcup_{i=1}^k S_i = G_0$. The next claim shows that our problem and the MSSC problem defined above are related.

Claim A.3. *Let $\{A_1, \dots, A_m\} \subseteq \mathcal{L}$ be such that $(\bigcap_{i=1}^m A_i) \cap G_0 = \emptyset$. Further, define $B_i = G_0 \setminus A_i$. Then the following is true*

$$\sum_{i=1}^t c_{\text{minsize}}(G_0 \cap A_1 \cap \dots \cap A_{i-1}, A_i) = \sum_{e \in G_0} \min\{i | e \in B_i\}. \quad (4)$$

Proof of Claim A.3. Note that by definition $|G_0| \leq |A_i|$ for every $1 \leq i \leq m$. Thus, for every $1 \leq i \leq t$, $|G_0 \cap A_1 \cap \dots \cap A_{i-1}| \leq |G_0| \leq |A_i|$. Thus, we have

$$\begin{aligned}
\sum_{i=1}^m c_{\min\text{size}}(G_0 \cap A_1 \cap \dots \cap A_{i-1}, A_i) &= \sum_{i=1}^m \min\{|G_0 \cap A_1 \cap \dots \cap A_{i-1}|, |A_i|\} \\
&= \sum_{i=1}^m |G_0 \cap A_1 \cap \dots \cap A_{i-1}| \\
&= \sum_{e \in G_0} \min\{i | e \notin A_i\} \\
&= \sum_{e \in G_0} \min\{i | e \in B_i\}.
\end{aligned}$$

□

Note that for any sequence A_1, \dots, A_m that satisfies the condition of Claim A.3, the sequence G_0, A_1, \dots, A_m is a valid solution to the list intersection problem on \mathcal{L} . Further, the cost of such a solution is the left hand side of (4). Also note that B_1, \dots, B_m is a valid solution for the MSSC problem on \mathcal{I} (as $\cup_{i=1}^m B_i = G_0$). The cost of this solution (for the MSSC problem on \mathcal{I}) is the right hand side of (4). Notice that L_1, \dots, L_m satisfies the condition of Claim A.3. Thus, by Claim A.3, this implies that there is a feasible solution for the MSSC problem on \mathcal{I} with cost OPT' .

Finally, consider the sets G_1, \dots, G_s . Note that the greedy algorithm for the MSSC problem will choose the lists from \mathcal{I} in the following order (for $1 \leq i \leq s$) $G_0 \setminus G_i$. Claim A.3 implies that $\text{GREEDY} = \text{GREEDY}'$, where GREEDY' is the cost incurred by $\text{GREEDY}_{\text{MSSC}}$. Thus, by Theorem A.1, Claim A.2 and the discussion above we have

$$\text{GREEDY} = \text{GREEDY}' \leq 4 \text{OPT}' \leq 8 \text{OPT},$$

as required. Thus, we shown the result for the special case when intersection of the lists in \mathcal{L} is empty.

Now assume that $I = \cap_{i=1}^k L_i \neq \emptyset$. For every $1 \leq i \leq k$, define $L'_i = L_i \setminus I$. Note that $\cap_{i=1}^k L'_i = \emptyset$. Further, notice that for any algorithm with inputs lists L_1, \dots, L_k , the cost paid at any step of the algorithm is composed of two parts. First, the algorithm pays a cost of $|I|$ and second, it pays a cost that depends only on the lists L'_i . Now, we decompose the total cost of the greedy and the optimal algorithm into two parts– GREEDY^1 , GREEDY^2 and OPT^1 and OPT^2 . Note that $\text{GREEDY}^1 = \text{OPT}^1 = (k-1)|I|$. Also by our analysis for the case when $I = \emptyset$, we have $\text{GREEDY}^2 \leq 8 \text{OPT}^2$. Summing up both parts of the costs incurred by the greedy and the optimal algorithm completes the proof of Theorem 2.1 in the general case.

We now turn to proving the hardness result in Theorem 2.1. We will reduce from the MSSC problem. Let the input instance to the MSSC problem be the lists S_1, S_2, \dots, S_m such that $U = \cup_{i=1}^m S_i$. Let \mathcal{S} denote the set of these lists. For every $1 \leq i \leq m$, let $T_i = U \setminus S_i$. Note that $\cap_{i=1}^m T_i = \emptyset$. Let U' be a set such that $U' \cap U = \emptyset$ and $|U'| = |U|$. Finally define the input lists for intersection problem as follows. $L_0 = U$ and for every $1 \leq i \leq m$, $L_i = T_i \cup U'$. Let $\mathcal{L}_{\mathcal{S}}$ denote these set of lists. Notice that $\cap_{j=0}^m L_j = \emptyset$.

Let $\text{OPT}(\mathcal{L}_{\mathcal{S}})$ and $\text{OPT}(\mathcal{S})$ denote the costs of the optimal solutions for the list intersection and MSSC problems respectively. We claim that

$$\text{OPT}(\mathcal{L}_{\mathcal{S}}) = \text{OPT}(\mathcal{S}) + |U|. \tag{5}$$

Theorem A.2 states that for the MSSC problem, it is NP-Hard to output a solution with cost at most $(4 - \varepsilon) \text{OPT}(\mathcal{S})$ for every input \mathcal{S} . Thus, it should be NP-hard to output a solution with cost $(4 - \varepsilon) \text{OPT}(\mathcal{S}) + |U|$ for the list intersection problem on $\mathcal{L}_{\mathcal{S}}$ for every \mathcal{S} . From (5), we have $\text{OPT}(\mathcal{L}_{\mathcal{S}}) = \text{OPT}(\mathcal{S}) + |U|$. Thus, it is NP-hard to approximate the list intersection problem to within a factor of

$$\frac{(4 - \varepsilon) \text{OPT}(\mathcal{S}) + |U|}{\text{OPT}(\mathcal{S}) + |U|} \geq \frac{5 - \varepsilon}{2},$$

where the inequality follows from the fact that $\text{OPT}(\mathcal{S}) \geq |U|$ (this follows from the definition of the MSSC problem).

To complete the proof, we show (5). We first note that the optimal sequence for the list intersection problem on \mathcal{L} must include L_0 (as $\cap_{i=1}^m L_i = U' \neq \emptyset = \cap_{i=0}^m L_i$). We now claim that the optimal solution must intersect L_0 (with some other list) in its first step. If not, w.l.o.g. let the sequence of lists considered be $L_1, \dots, L_s, L_0, L_{s+1}, \dots, L_m$. Instead consider the sequence $L_0, L_1, \dots, L_s, L_{s+1}, \dots, L_m$. Note that after s steps, both the sequences incur the same cost. Note that $U' \subseteq \cap_{i=1}^s L_i$. Thus, in the first sequence, in each of the first s steps, the sequence pays at least $|U'| = |U|$. The second sequence in the first step pays $\min\{|L_0|, |L_1|\} = |U|$. Now note that $(L_0 \cap L_1) \subseteq U$ and thus, in all of the remaining steps, the second sequence pays at most $|U|$. Thus, the second sequence has a total cost no worse than the first sequence.

Thus, we now have that the optimal solution for intersecting lists from \mathcal{L} intersects L_0 in its first step. In other words, after the first step, the algorithm is solving the list intersection problems on the sets T_1, T_2, \dots, T_m . By Claim A.3, $\text{OPT}(\mathcal{T}) = \text{OPT}(\mathcal{S})$, where \mathcal{T} is the list intersection problem instance on the sets T_1, \dots, T_m . This along with the fact that $\text{OPT}(\mathcal{L}_{\mathcal{S}}) = |U| + \text{OPT}(\mathcal{T})$ proves (5). \square

B Proofs: Comparisons Model

Proof of Theorem 4.1. Let the lists in the input \mathcal{L} be denoted by L_1, L_2, \dots, L_k . We will assume that the optimal AND-tree is left-deep (we will see later how to get rid of this assumption). We first look at the case when the intersection is empty, that is, $\cap_{i=1}^k L_i = \emptyset$.

Let G_0, G_1, \dots, G_s be the lists used by the greedy algorithm. Note that by the definition of the greedy algorithm, $|G_0| = \ell_{\min}$ (and also by the assumption on \mathcal{L} , $\cap_{i=0}^s G_i = \emptyset$). Further, let GREEDY_j be the cost incurred by the greedy algorithm on step $j = 1, 2, \dots, s$. We do not have a precise handle on the number of comparisons made at step j . However, notice that we can upper bound it by twice the size of the smaller list. Thus, for $1 \leq j \leq s$,

$$\begin{aligned} \text{GREEDY}_j &\leq 2 \min\{|G_0 \cap G_1 \cap \dots \cap G_{j-1}|, |G_j|\} \\ &= 2|G_0 \cap G_1 \cap \dots \cap G_{j-1}| \end{aligned} \tag{6}$$

$$\leq 2\ell_{\min}, \tag{7}$$

where in the above we have used the facts that $|G_0 \cap G_1 \cap \dots \cap G_{j-1}| \leq |G_0| \leq |G_j|$.

Now let us look at the optimal sequence. W.l.o.g. assume that the sequence is L_1, L_2, \dots, L_t . For step $1 \leq j \leq t - 1$, let OPT_j denote the cost paid by the optimal solution in that step. Again, we do not have a precise handle on the number of comparisons made by the optimal algorithm at step j . However, note that we can lower bound the number of comparisons made by the size of the

resulting intersection at the end of step j . Thus, we have

$$\begin{aligned} \text{OPT}_j &\geq |L_1 \cap L_2 \cap \dots \cap L_{j+1}| \\ &\geq |G_0 \cap L_1 \cap L_2 \cap \dots \cap L_{j+1}| \end{aligned} \quad (8)$$

Note that now we are in the familiar terrain of the c_{minsize} function (we lose a factor of 2 in the process). As in the proof of Theorem 2.1, we consider the following instance of the MSSC problem. The universe is G_0 and the lists are defined as $S_i = G_0 \setminus L_i$. Call this instance \mathcal{I} .

First we claim that for $j = 1, \dots, s$ the $\text{GREEDY}_{\text{MSSC}}$ algorithm chooses the list $G_0 \setminus G_j$ in step j . This is because by definition of the greedy algorithm for the list intersection problem, G_j has the smallest intersection (among lists in $\mathcal{L} \setminus \{G_0, \dots, G_{j-1}\}$) with $G_0 \cap G_1 \cap \dots \cap G_{j-1}$. In other words, $G_0 \setminus G_j$ covers the maximum number of elements in $G_0 \setminus (G_0 \setminus G_1 \cup \dots \cup G_0 \setminus G_{j-1})$. Thus, $\text{GREEDY}_{\text{MSSC}}$ will pick $G_0 \setminus G_j$ in step j , as claimed.

Let GREEDY' be the cost incurred by the $\text{GREEDY}_{\text{MSSC}}$ sequence on \mathcal{I} . Thus, we have

$$\begin{aligned} \text{GREEDY}' &= \sum_{e \in G_0} \min\{j | e \in G_0 \setminus G_j\} \\ &= |G_0| + \sum_{j=1}^s |G_0 \cap G_1 \cap \dots \cap G_j|, \end{aligned} \quad (9)$$

where the second equality follows from the following accounting. GREEDY' is the sum over all elements in G_0 the first step in which the element is covered (this is the first equality above). The index of the first step is the same as one plus the number of steps for which the element is *not* covered. Summing this up over all elements and rearranging the summations gives the second equality. For notational convenience define $\text{GREEDY}'_j = |G_0 \cap G_1 \cap \dots \cap G_j|$. By (6) we get for all $2 \leq j \leq s$,

$$\text{GREEDY}_j \leq 2 \text{GREEDY}'_j. \quad (10)$$

Now consider the optimal sequence L_1, \dots, L_t with cost OPT . By arguments similar to the ones used to relate the greedy algorithms, the sequence $G_0 \setminus L_1, \dots, G_0 \setminus L_t$ is a feasible solution to the MSSC problem on \mathcal{I} . Let OPT' denote the cost of this solution. Again using arguments similar to the ones used to derive (9), we get

$$\text{OPT}' = |G_0| + \sum_{j=1}^t |G_0 \cap L_1 \cap \dots \cap L_j|. \quad (11)$$

Again for notational convenience define $\text{OPT}'_j = |G_0 \cap L_1 \cap \dots \cap L_j|$. By (8), we get that for every $1 \leq j \leq t-1$,

$$\text{OPT}_j \geq \text{OPT}'_{j+1}. \quad (12)$$

By Theorem A.1, (9) and (11) we have the following inequality.

$$\ell_{\min} + \sum_{j=1}^s \text{GREEDY}'_j \leq 4 \left(\ell_{\min} + \sum_{j=1}^t \text{OPT}'_j \right). \quad (13)$$

Due to lack of space, we omit the proof for the case when the intersection is empty.

The following sequence completes the proof for the case when the intersection is empty.

$$\begin{aligned} \text{GREEDY} &= \sum_{j=1}^s \text{GREEDY}_j \\ &\leq 2\ell_{\min} + \sum_{j=2}^s \text{GREEDY}_j \end{aligned} \tag{14}$$

$$\leq 2\left(\ell_{\min} + \sum_{j=1}^s \text{GREEDY}'_j\right) \tag{15}$$

$$\leq 8\left(\ell_{\min} + \sum_{j=1}^t \text{OPT}'_j\right) \tag{16}$$

$$\begin{aligned} &= 8\left(\ell_{\min} + \text{OPT}'_1 + \sum_{j=2}^t \text{OPT}'_j\right) \\ &\leq 8\left(\ell_{\min} + \ell_{\min} + \sum_{j=1}^t \text{OPT}_j\right) \end{aligned} \tag{17}$$

$$= 8 \text{OPT} + 16\ell_{\min}.$$

(14), (15) and (16) follow from (7), (10) and (13) respectively. Finally, (17) follows from (12) and the fact that for every j , OPT'_j is upper bounded by the size of the universe, which is G_0 .

To complete the proof we need to consider the case when $I = \cap_{i=1}^k L_i \neq \emptyset$. As in the proof of Theorem 2.1, the argument for the general case (given the analysis for the case $I = \emptyset$) divides up the costs for both the greedy and optimal algorithms as GREEDY^1 , GREEDY^2 and OPT^1 , OPT^2 . Recall that GREEDY^1 and OPT^1 depend only on I . Again using arguments similar to those used to derive (6) and (8), we get the bounds $\text{GREEDY}^1 \leq 2(k-1)|I|$ and $\text{OPT}^1 \geq (k-1)|I|$. By our analysis for the case when $I = \emptyset$, we get $\text{GREEDY}^2 \leq 8\text{OPT}^2 + 16\ell_{\min}$. Summing up bounds on both parts of costs incurred by the greedy and the optimal algorithm completes the proof for the general intersection case.

We now briefly argue why we could assume that the optimal AND-tree was left-deep. For the comparison model, it is not clear why the optimal AND-tree has to be left-deep. However, note that our proof actually proves that the greedy algorithm in the comparison model is within a constant factor of the optimal AND-tree in the intersection size model, that is, for an intersection, we charged the optimal algorithm just the size of the intersection. Under the modified cost model an argument very similar to that in the proof of Claim A.1, shows that the optimal AND-tree is a left-deep tree. The crux of the argument is to show the result for the following special case. Let the optimal AND-tree correspond to first intersect list A and B , then intersect C and D and finally intersect the two intermediate lists. W.l.o.g. assume that $|A \cap B| \leq |C \cap D|$ and that A and B are in the left sub-tree at the root. Now consider the following ordering: first intersect A and B , then intersect the result with C and finally intersect the intermediate list with D . It is easy to check that the latter is a left-deep tree and its total cost is at most that of the original tree.

We now turn to the proof of the lower bound, which is constructive. Let $\Sigma = \{0, 1\}^k$. Any element $x \in \Sigma$ will be associated with a vector (x_1, \dots, x_k) where every $x_i \in \{0, 1\}$. We define the lists as follows. For every $x \in \Sigma$, $x \in L_1$ ($x \in L_2$) if and only if $x_1 = 0$ ($x_1 = 1$). For any

$i = 3, \dots, k$, for every $x \in \Sigma$, $x \in L_i$ if and only if $x_i = 0$.

First note that $L_1 \cap L_2 = \emptyset$. Further, no other collection of lists leads to \emptyset . Thus, the optimal solution has to intersect these two lists— further, it is easy to see that the optimal solution can do it with *one* comparison.

Now, it is not too hard to see that for every i , $r \geq 1$ and set of indices $I = \{i_1, i_2, \dots, i_r\}$ such that $|I \cap \{1, 2\}| \leq 1$, the following holds: $|\cap_{j \in I} L_j| = 2^{k-r}$. Thus, for every (but the last two) step, all the lists are “equivalent” for the greedy algorithm. Thus, $L_3, \dots, L_k, L_1, L_2$ is a valid sequence in which the greedy algorithm intersects the lists. Thus, after the $1 \leq i \leq k - 3$ intersections, the size of “current” list will be $2^{k-i-1} = 2^{-i} \cdot \ell_{min}$. Now for each intersection, the greedy algorithm has to make at least as many comparison as the size of the final intersection. Thus, the number of comparisons made by the greedy algorithm is at least

$$\frac{\ell_{min}}{2} \left(1 + \frac{1}{2} + \dots + \left(\frac{1}{2} \right)^{k-4} \right) = \ell_{min} \left(1 - \left(\frac{1}{2} \right)^{k-3} \right).$$

□