

VDM Aggregation Strategy

By Michael Kamfonas

A successful performance strategy for very large databases (VLDB) often involves aggregations. This is why major DBMS vendors offer mature system support for aggregation maintenance and optimization. DB2 offers Materialized Query Tables (MQT), Teradata offers Aggregate Join Indexes (AJI), Oracle offers Materialized Views (MV). BI tool vendors support aggregate awareness and dynamic sourcing, while OLAP technologies offer cost-effective, in-memory retention alternatives. All too often, however, these powerful capabilities are used redundantly, without too much forethought, and although they often improve query performance, they sometimes also disappoint, by impacting ETL windows, availability, stability, sustainability and sometimes quality. This is an overview of InfoKarta's approach to implementing an Aggregation Strategy. It proposes a model and a framework to reason about aggregates, qualify work-loads that benefit from them, and pick most effective levels from a total performance perspective. It also recommends design and implementation approaches for exploiting the DBMS aggregation technology at hand in a way that is resilient, sustainable, scalable and maintainable. The techniques described are part of the Versioned Dimensional Model¹ (VDM) Toolkit and methodology which originated in the early 90s by the author, and has been evolving ever since.

Why an Aggregation Strategy?

The conversation usually goes something like this:

Me: We need to optimize for Sales Dashboard queries at any level of Location, Product and Time Period. We should be able to support minimum service levels across the envelope of all drill-down summaries that these queries can hit.

DBA: Even very similar queries may follow different execution plans. We need to optimize for the levels and types that are most critical. Let's get a list of the 10-20 typical or worst offender queries to optimize. Then we will clearly be able to measure success and know when we are done. Then we can repeat the exercise.

Me: Even if we pick 20 queries, we know that the patterns will change or new ones will emerge. We need to ensure that the whole envelope of dashboard queries has good response. All these queries are one or two-level drilldowns, and must consume less than 150 CPU-seconds average to sustain the required concurrency. At the same

¹ The Versioned Dimensional Model (VDM) comprises methods and techniques around the design and implementation of analytic and reporting databases, bridging the gap between operational and dimensional models with emphasis on performance on parallel share nothing platforms. The VDM-Toolkit© is a framework of customizable tools and accelerators around the author's VDM practice.

time we have to make sure that heavily hit levels are supported with close enough aggregates to reduce total cost.

DBA: This is too open-ended. Query patterns are too nebulous. We cannot optimize for queries we can't see or run. I say, find the most common queries and optimize for those. We can continue monitoring the workload, find new worst offenders the following week and optimize them in the next round. Iterative improvement is the name of the game!

This article was conceived to articulate how optimizing for aggregation envelopes and query patterns can be conceived, approached systematically, measured and architected.

If this thought appeals to you, or any of the situations below sound familiar, you should read-on:

- Overnight maintenance windows get tighter and tighter. Opening times start getting missed as more and more aggregations elongate ETL cycles and squeezing contingency margins.
- Unexpected delays of alarming magnitude started to occur occasionally. Aggregation maintenance can add unpredictability to ETL time windows. Small dimension changes, for example, can cause disproportionately large and expensive updates when recasting is involved.
- Inconsistencies between base and aggregates cast doubt in users' mind about the quality of data. Missed or delayed maintenance causes discrepancies to become visible through reports or dashboards. Adding insult to injury, it is often the users that discover these issues and tracing the root cause can be time-consuming and sometimes inconclusive.
- Newly introduced reports or query patterns, with no aggregate support, abuse resources saturating the system and impacting other production workloads. Existing aggregate JIs or MQTs need to be adjusted or new ones introduced.
- After two years of active use, it is not uncommon for a data warehouse to lose a third or more of its capacity in redundant or unnecessary copies of data. As migrations to newer versions of base or derived data occur, some applications remain coupled to older, now redundant, data objects. Application change is required, hampering the elimination of old data which as time goes by and the focus of key developers moves to other areas it becomes riskier to do. Inevitably, the only viable solution is to upgrade capacity earlier than planned. The reasons for data redundancy are many, and aggregate-to-application coupling is one of them.
- OLAP and in-memory cube technologies reduce the costs of high concurrent usage and can be welcome additions to our arsenal. But how does one prevent rules and metric definitions from drifting into inconsistent parochial application meta-models and enforce consistency across all access tools in the data warehouse environment?

The Problem Space

Data warehouses, data marts and BI tools enable a chain of model transformations that start from the source normalized model, and end in the myriad of result sets delivered to end-applications and users. The architect has to decide which “snapshots” of this transformation chain to instantiate into concrete materialized physical objects, effectively dividing the chain into daily batch and on demand segments. Figure 1 shows this continuum and how the options explode as we move away from the source model and towards the right, the consumption-end of the warehouse. The most popular choice is to materialize a base model close to the source model. This not only favors ETL simplicity but it also preserves, relatively untampered, the original content satisfying the broadest spectrum of requirements either not yet uncovered or needing correction or reinterpretation. Downstream transformation from the base model ideally could be dynamic, through one or more layers of views, delivering an access layer organized by functional area of interest. This access layer often presents a dimensional model that makes it easy for applications and users not only to select and filter, but also to aggregate and combine facts to derive complex metrics. Keeping this transformation dynamic is elegant, maximizes availability and keeps the pain of change low since the base model is relatively normalized. However, even with ideal indexing and physical organization, there is too much transformation left in the pipeline to absorb at query time, leading to performance degradation, saturation of resources and excessive cost.

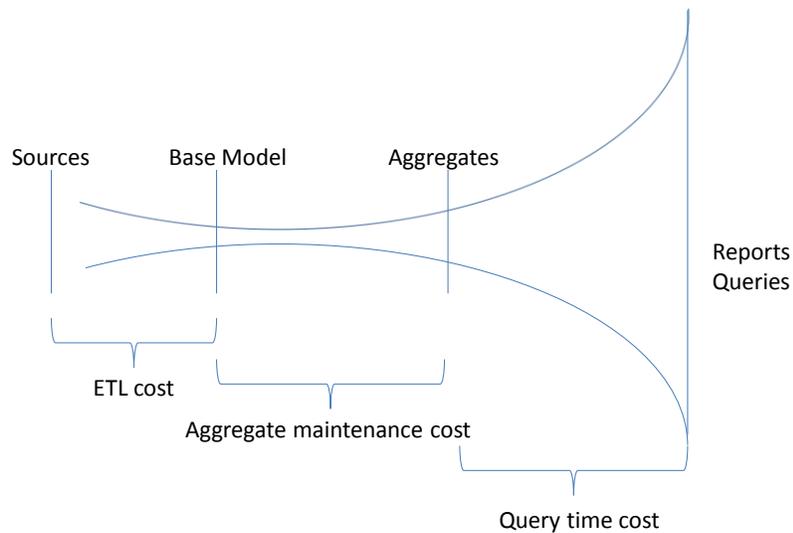


Figure 1 Transformation pipeline and choice of materialization points

To mitigate the problem we seek appropriate intermediate materialization points along the transformation path to take some of this burden away from the query, process it in advance, typically once a day, and share it across multiple more efficient query executions. The improvement most typically comes by reducing I/O in the form of pre-joins and aggregations. The further to the right of the diagram these aggregates fall i.e. the closer to the targeted reports, the more efficient queries will become; but also the more explosive the encountered patterns, and thus the cost of aggregate maintenance. The VDM Aggregation Strategy is a methodology that finds and implements aggregates to optimize holistic system performance. It is appropriate where a large portion of the query workload aggregates data and follows a dimensional pattern, even if the schema is not dimensional.

The essence behind a dimensional query pattern is in the aggregation of lower level events along the apparently independent (but potentially correlated) classifiers and taxonomies they are joined with. The schema best suited for this purpose is highly dependent on the DBMS support in terms of optimization capability and implemented access methods. Typically, the CPU and I/O costs are highly correlated to the number of rows scanned. Our optimization approach seeks to minimize row counts by pre-aggregating data, while also combining other pragmatic considerations to arrive at a rational, flexible solution.

The three obvious questions, when it comes to aggregations, are:

- What are the optimal levels?
- How should we implement and maintain them?
- How can we exploit them so our queries run predictably faster?

In support of the last two bullets, major DBMSs (DB2, Teradata and Oracle) offer methods for defining and maintaining aggregates automatically, and provide some type of re-write capability during optimization to redirect qualifying queries to the appropriate aggregate. There are nuances in the implementations and each DBMS vendor puts forth good reasons on why they do or don't allow certain features. For example, Teradata only allows system-maintained Aggregate Join Indexes (AJIs) keeping the implementation simpler and always ensuring consistency between base tables and aggregate. IBM/UDB and Oracle allow more flexibility, adding options for deferred or user-maintained MQTs and MVs. This flexibility comes at the cost of more complexity and risk exposure. Of course, good old-fashioned aggregate tables with no DBMS support are always available and OLAP, in-memory cube technologies or BI-tool capabilities add to the menu of choices.

The successful aggregation strategy is less about features, and more about how well we exploit the capabilities available to us. The approach described below combines analytical optimization techniques with technology-specific practices, and pragmatic considerations for complexity, stability and sustainability.

The VDM Aggregation Strategy Framework

The “Aggregation Lattice”

We use two types of graphic representations to aid us in visualizing aggregations. The “Level Grid” illustrated in Figure 2 shows a fact (in green) with additive metrics M1, M2, against the five dimension keys D1-D5. The dimensions are represented by vertical lines stemming off these keys with their levels numbered starting from zero.

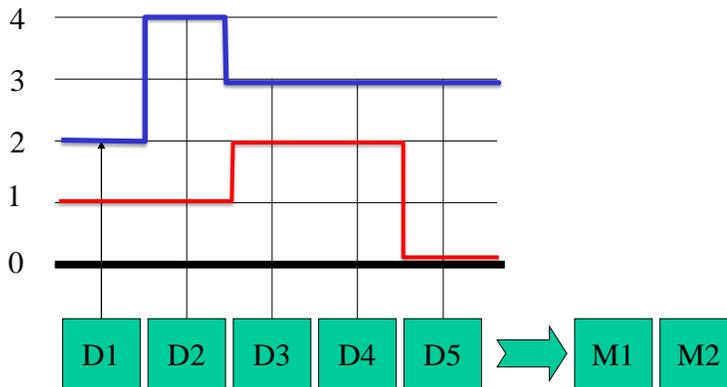


Figure 2 - Basic Dimensionality of our example fact

The black line at the bottom denotes the base level of the fact, while the blue line at (2,4,3,3,3) indicates the top aggregate along all dimensions – a single row summary. Any line which cuts across the five dimensions at a valid level represents an aggregation. The red one shown in the Figure 2 is at level (1,1,2,2,0).

The “Aggregation Lattice” is the second type of graph we use, depicting all possible aggregations from the base fact to the highest level. Every aggregation-defining line that cuts across our “level Grid” corresponds to a node of the Aggregation Lattice. The lattice also has edges that indicate all possible derivation paths of each aggregate from lower levels. Our example in figure 1 has $3 \times 5 \times 4 \times 4 \times 4 = 960$ possible such lines of dimension level combinations or 960 nodes in the corresponding Aggregation Lattice. Since such size lattice would be too cluttered to work and visualize, Figures 3 and 4 show smaller examples with two and three dimensions respectively. Here are the conventions we use to construct these lattices:

1. Each dimension is a tree of uniform depth, with its nodes organized in levels, with a single top node representing “all” or “any”. Leaf nodes are at level 0; their parents are at level 1; their parents at level 2 etc. The root (top) node level is $n - 1$, where n represents the number of levels in the dimension. Simple codes with no hierarchical structure can be represented in two levels: the code itself at level-0, and a single parent denoting the “any code” at level 1.
2. There are as many possible aggregates as there are combinations of participating dimension levels. We use the term “multi-levels” when we want to distinguish them from single dimension levels. A bottom multi-level for a three-dimensional model is $(0,0,0)^2$ and it represents the base fact; the top can be something like $(4,5,4)$ representing all stores, all product, for all time. All possible multi-levels between top and bottom are connected by arcs depicting every possible single-dimension single-level aggregation path. The graph so formed is a lattice, “the Aggregation Lattice,” a complete map of all valid aggregation paths.

Three lattice examples of a two-dimensional hierarchy are depicted in Figure 3. The number of nodes (vertices for the mathematically inclined) is the product of the number of levels for each dimension; from left to right, $2 \times 2 = 4$, $4 \times 4 = 16$ and $3 \times 3 = 9$. Similarly, Figure 4 depicts three-dimensional lattices for two and three level hierarchies with $2 \times 2 \times 2 = 8$ and $3 \times 3 \times 3 = 27$ nodes respectively. For illustration purposes we show transitive aggregations using dashed lines for the 3×3 lattice.

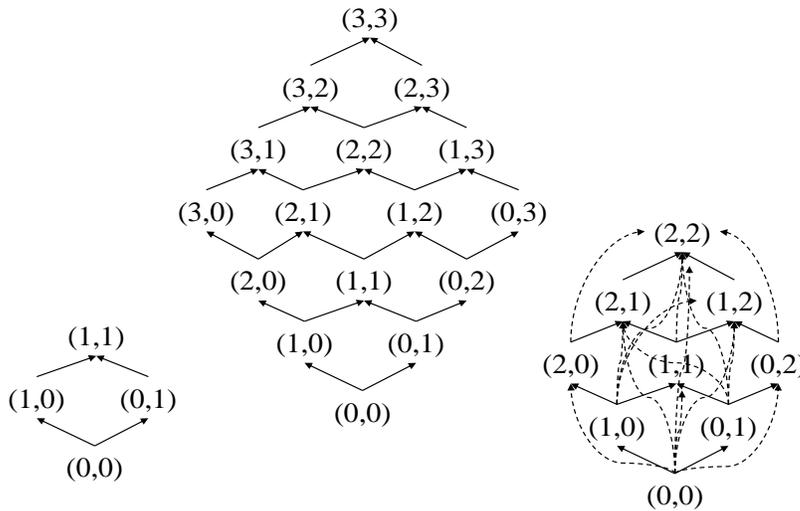


Figure 3: Two-dimensional aggregation lattices 2, 3 and 4 levels deep.

² There is a subtle difference between the base fact $(0,0,0)$ and the transactional detail level, where all event data are stored. The methodology uses an additional node, referred to as level -1 below the $(0,0,0)$ level to denote the non-aggregated underlying event data. For simplicity, we ignore it during this discourse.

There is more than one way to derive an aggregation from lower levels, and the number of ways increases as the number of dimensions increase. For example, (2,0,2) is derived by aggregating one level along the third dimension of (2,0,1) or along the first dimension of (1,0,2), or by aggregating any of the lower levels, i.e. (2,0,0), (1,0,1), (0,0,2) or (1,0,0), (0,0,1) and (0,0,0).

In the diagrams we show vertices in layers, each representing multi-levels of the same “rank” or “aggregation distance” from the bottom. The lowest, single vertex represents the base fact and has rank 0; the next layer-up has vertices with rank 1, and so on. The rank correlates to the “level of magnitude” of an aggregate in terms of row count. To illustrate this, visualize an ideal mart where all dimensions are independent with same depth and cardinality ratio 10:1 from level to level. If the base fact has 1 million rows the vertices in the next layer represent multi-levels (or aggregates) with rank 1 and 100,000 rows. The next layer of rank 2 will have cardinalities ten times less and so on. Of course in reality dimensions are not independent and cardinality ratios vary. This causes aggregates of the same rank to have very different row-counts. This is why we use the notion of “rank” to talk about the topology, but we compute actual cardinalities to analyze performance potential.

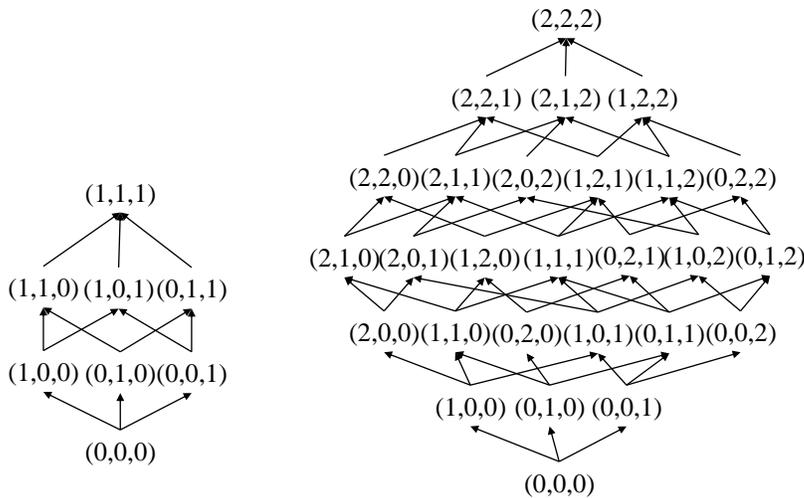


Figure 4 Tree-dimensional aggregation lattice 2 and 3 levels deep.

The first step of our process is to form a “Loaded, Weighed Aggregation Lattice” as follows:

1. Establish the dimensionality of each fact. This is a critical step, particularly if the underlying schema is non-dimensional. This entails organizing the model into dimensions, picking the relevant levels and pruning those dimensions that we should not consider or that have no value for aggregate construction.
2. Formulate “aggregation lattice” of all level combinations along the dimensions.
3. Load nodes with cardinalities derived from actual data sampling to arrive at the “Loaded Aggregation Lattice”
4. Consider additional expected or projected usage, physical organization constraints etc., arriving at one or more “Weighted Loaded Aggregation Lattices” which in turn are used to select optimal aggregate candidates.

Query Patterns and Aggregation

In order for the aggregation lattice to be useful, we need to credibly link it to query performance. What query patterns, and what types of workloads are conducive to performance improvement through aggregates?

Every query, however complex, is associated with four multi-levels as shown in figure 4:

1. **The Selection/filtering multi-level**, is the lowest applied filtering level along each of the dimensions. This is determined by the “where” clause of an SQL query. For dimensions that don’t appear in the where-clause, we assume filtering at the top level of the hierarchy representing “all” or “any” is implied. E.g. filtering for product category A for 2015Q1, automatically assumes “all stores”. We refer to this multi-level as the S-Level (for Selection Level)
2. **The Grouping multi-level** is the lowest level along each dimension found in the “group by” clause of the query. We refer to this level as the G-Level.
3. **The Materialized Table multi-level** is the actual level of the table being queried – the M-Level (Materialized Data level.) The materialized level, in the absence of aggregates, is the base fact level. The further apart S-Level and M-Levels are, the higher the cost of the query. Intuitively, the higher the filtering level means more rows to scan; the lower the materialized level the more granular the data we have to scan. Figure 5 shows our fact with five dimensions, the S-Level (dotted), G-Level (solid) and the M-Level, in this case (0,0,0,0,0).

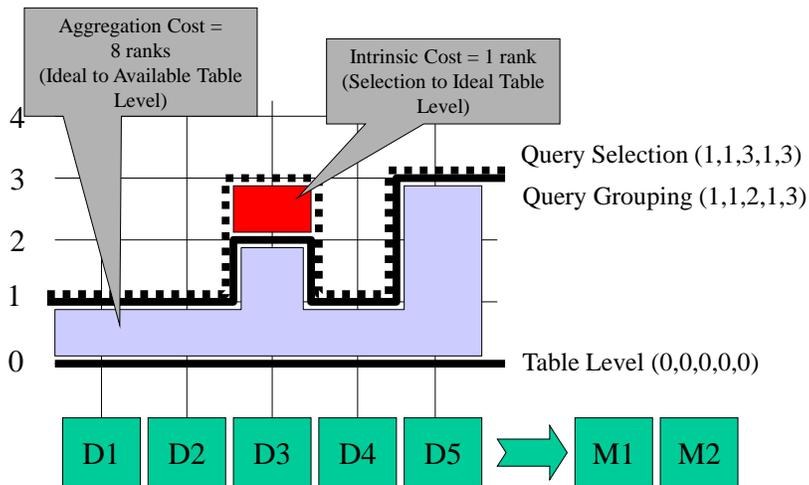


Figure 5 Query profile over a 5-dimensional aggregation hierarchy.

4. **The Query multi-level** is the ideal materialized data aggregate the query could run on (Q-Level), if it existed. It usually coincides with the G-Level but not always. The query level is formed by checking each dimension S and G levels and picking the lower of the two.

Figure 6 depicts the same scenario shown in Figure 5 with an aggregate added at level (1,1,1,1,2). The **Total Aggregation Cost** from base data to the query result corresponds to the area between S-Level and Base and it ties back to the sum of Aggregate Maintenance plus Query Run-Time Cost of the transformation chain shown in Figure 1. We breakdown this aggregation cost into component costs:

- **Pre-Aggregation Cost:** The cost to pre-aggregate data from the base level to the available M-Level aggregate used by the query, i.e. between the S-Level and M-Level. This corresponds to the Aggregate Maintenance Cost in Figure 1.
- **Run-Time Aggregation Cost:** This cost is represented by the area between the S-Level and the M-Level, where data actually is queried from. It corresponds to the Run-Time Aggregation Cost in Figure 1. The Q-Level divides this area into two parts, thus further breaking this cost down to two component costs:
 - **Intrinsic Aggregation Cost:** The portion of the aggregation that is inherent to the query and no aggregate can reduce. It is represented by the area between the S-Level and the Q-Level. Low intrinsic cost characterizes queries whose scan size can be relatively small if a good aggregate can be provided for them to run on. Conversely, high intrinsic

- cost queries require large numbers of qualifying rows scanned per point selected, even if the best possible aggregate existed for them to run on.
- **Discretionary Aggregation Cost:** This is the run-time cost incurred to aggregate data from an available materialized aggregate up to the query level, i.e. the area between the Q-Level and the M-Level. It is discretionary since we chose to materialize at a lower level. Essentially, introducing aggregates converts Discretionary to Pre-Aggregation cost.

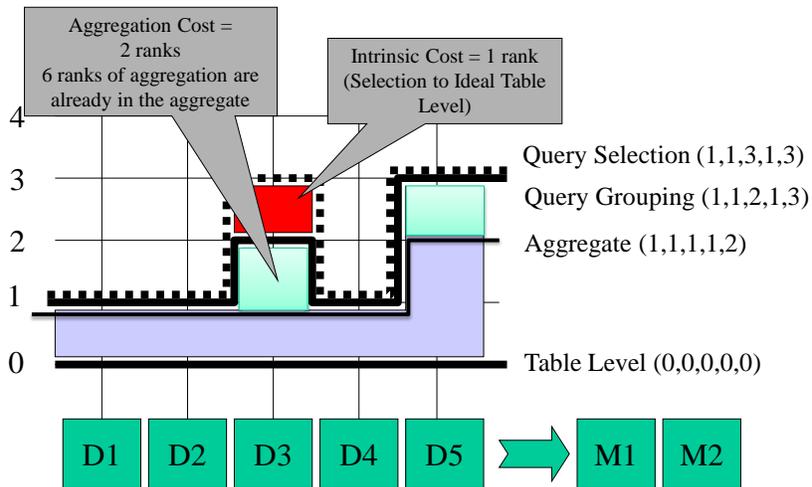


Figure 6 - Intrinsic, discretionary aggregation and pre-aggregation cost separated

We can classify aggregation queries into distinct patterns based on the relative positioning of these lines and their intrinsic cost. Figure 7 shows four such patterns:

1. Drill-Down Interactive Queries are typically initiated from a point the user selects and requests a lower level drill down. The selected cells are at some “selection level” (S-Level) and the drill-down level defines the “grouping level” requested (G-Level.) The drill-down is typically the next level along one or two dimensions. In lattice terms, the S and G levels are close to each other, maybe a couple of ranks apart, meaning that only a few G-Level rows are required per S-Level selected. In other words, the intrinsic cost is low and consequently sustaining interactive response times is feasible as long as an aggregate exists “close enough” to any query level selected. The top left diagram of the figure shows how from a selection level at (3,2,4,3,1) drilling down along two dimensions delivers grouped data at (3,2,3,2,1). The red shaded area is proportional to the intrinsic cost of the query.

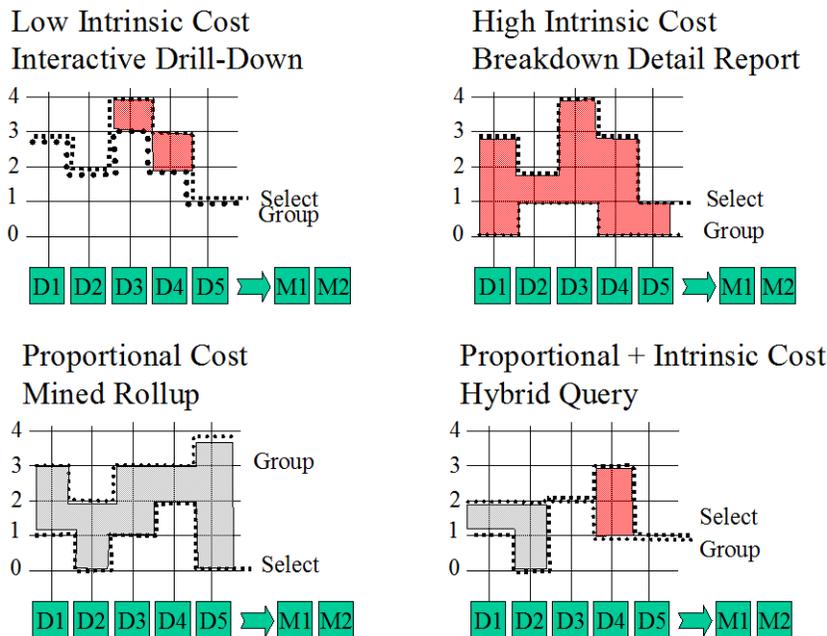


Figure 7: Query patterns

2. Breakdown Reports are characterized by a relatively broad selection and relatively low-level grouping of data e.g. select Product Categories X, Y and Z for last month for this and last year, and return amounts and quantities sold by product class, by store and day. Such a pattern is shown at the top-right diagram of Figure 7. These queries have a high intrinsic cost because of the large distance between selection and grouping levels. They can only be run on relatively low level detail and they require a large number of rows. They are classified as “reports” since they can be rather voluminous due to the level of detail they convey. This type of query may be used to feed graphic visual components to plot highly detailed graphs, in those cases naming their classification as “reports” may be too restrictive.
3. Mining and Hybrid Rollups are queries that select data along some or all dimensions at levels lower that they group them. The lower diagrams of Figure 7 show such query patterns mining low level data and rolling them up to cluster or classify them along higher levels of the same dimensions. Select specific SKUs supplied by a list of vendors last year and group them by class and district. This type of query needs to mine data for specific products at the item level, and then roll-up counts and amounts to higher levels of these taxonomies.

With this backdrop, let’s now look into how optimal levels are chosen and the details behind their implementation.

Optimal Levels

Formality and General Practice

One school of thought is to “architect” aggregates during the data design process and base them on application/reporting requirements that dictate what the hot spots are. This inevitably results in hardwired aggregates, often modeled as prime structures and taking life of their own, difficult to adjust and likely to leave large portions of the lattice uncovered.

A different approach is to monitor the system workload and create aggregates when needed in response to performance issues when they arise. This approach is adaptive and responds to changing workloads, however it may lead to proliferation of costly high proximity application-specific aggregates that are easy to add but hard to refactor. This in turn is likely to interfere with their sustainability and evolution, leading to resource saturation and premature need for capacity upgrades.

Analytic Approach

The approach we favor combines the “architected” with the “iterative” and it is based on the following three premises:

- Optimal aggregate selection is driven from the data itself; not directly dictated by specific reporting requirements.
- A rational framework for quantifying and analyzing performance is essential, and optimization should take into account the total system, its workloads and service levels.
- Coupling of applications to aggregates must be avoided so that changes are painless down the road.

We use two types of algorithms to determine optimal aggregates. One is a variation on the technique proposed by Harinarayan, Rajaraman, and Ullman [1], which iteratively seeks the “best next aggregate” that minimizes the sum of materialized rows scanned to run the equivalent of “select *” at every vertex of the lattice. As higher level aggregates get picked and their materialization factored into the model, this sum is reduced progressively. The other is a VDM technique that sweeps the weighted loaded aggregation lattice starting from the bottom, to select the closest node of “adequate aggregation distance.” The algorithm iterates and as chosen aggregates are factored into the model all vertices of the lattice end up “close enough” to some chosen aggregate thus covering the whole lattice with heavier weighed nodes favoring hot spots.

The purpose of this paper, however, is less to explain algorithms, and rather to present and rationalize the methodology, its premises and characteristics:

1. Optimal aggregations are best determined by the data topology and cardinalities depicted in a properly weighed and loaded aggregation lattice. Requirements should

be used to construct, prune, load or weigh the lattice, but should not dictate aggregate levels.

2. Maximize sharing. One common aggregate is better than two close ones over the same fact supporting two different applications. There is one caveat to this, having to do with autonomy and separation of concerns, and it is discussed separately in the next section.
3. Odd but true: As a general rule the more effective the aggregates chosen to materialize, the more expensive they will be to maintain. Chains of aggregates built one on top of the other facilitate maintenance, but tend to leave large areas of the lattice uncovered. Proper aggregate selection should provide complete coverage of the lattice and address hot spots by weighing them accordingly.
4. Avoid trivial aggregations. Aggregating 100 billion rows into a 90 billion row aggregate, at face value, makes a negligible improvement in performance and is not worth the cost of its maintenance. Such close aggregates may only be justified if they resolve other costly derivations, such as:
 - a. Pre-joining orders to Line items
 - b. Pivoting vertically organized health provider roles into flat easier queried columns
 - c. Materializing derived columns that are used for filtering.

Such situations can often be addressed in the base model design, but by the time they become an issue aggregates are the easier way to resolve them.

5. Avoid recastable aggregations if possible. If an as-is dimension participates in the lattice, the lowest and highest levels are by definition non-recastable. Levels in between may require restating historical aggregates if the hierarchy changes. Evaluate the feasibility of pruning recastable multi-levels from the aggregation lattice. Although good to avoid, recastable aggregates may be a necessary evil to achieve desired service levels, and they should not be discarded a priori.
6. Follow the rule of proportionality: The magnitude of the work executed by the system to ingest a data change and maintain all dependent structures must be proportional to the magnitude of the change ingested. This is a generalization of the earlier point about recasting. Examples where this rule may get violated are:
 - a. A change of one row of a dimension may impact millions of fact rows (as-is recasting)
 - b. The use of MIN or MAX in an aggregate may cause history refresh if a row that determined the minimum or maximum is removed
 - c. Data maintenance practices - Refreshing a dimension by replacing data as opposed to merging changes will cause the whole dependent tower of dependent JIs to be refreshed.

In most cases, proportionality violations can be fixed by tweaking scripts, adding an index, running statistics, setting up proper constraints etc. It is important that these issues get shaken out of the system before aggregates are considered. Aggregates should not be used as a crutch or workaround to basic performance tuning.

Implementation Considerations

Aggregate Maintenance Automation – The Choices

There is no doubt that system-maintained data is the preferred method from the point of view of data integrity and quality. From the perspective of performance, however, it can be argued that user-maintenance methods can outperform automated methods in some cases. For example, system maintenance that is row-at-a-time or constrained by transaction boundaries leads to fragmentation, additional overhead and the possibility of double updates where independent details fall under the same summary point. Deferred system maintenance, if supported, can be an improvement. We encourage systematic testing and relentless tuning and qualification of the available system-provided methods before resorting to user-built solutions.

As a rule we discourage user-controlled solutions when system-controlled options are available. There are cases, however, where materializing actual tables is appropriate. The pros and cons vary by platform:

- If an aggregate involves metrics that violate the proportionality principle, such as MIN and MAX, it will not be incrementally maintained by the system, resorting to full refresh.
 - In DB2 a user maintained MQT is preferable because it can still be recognized for optimization.
 - In Teradata a user-maintained table may be a better choice with BI tools providing query optimization through aggregate awareness features.
- If a recastable (as-is) dimension is involved in an aggregation and we want to avoid the unpredictability of real-time maintenance caused by dimension changes.
 - DB2 deferred MQTs may provide adequate flexibility. User-maintained MQTs may be a better and more predictable option. A VDM differential recasting algorithm is provided.
 - In Teradata deferred maintenance of JIs is not currently supported. If the JI exposes unpredictability risk, a user-maintained aggregate table should be considered, relying on BI tools for optimization.
- The complexity of queries, after views are unraveled, may impact optimization resulting in suboptimal plan selection, reduced estimation confidence or failure to match JIs and MQTs. Low level facts can be materialized as independent tables absorbing computation and join complexity. Simpler aggregates can then be built on top.

Sharing, Autonomy and Separation of Concerns

Two data marts that share a sales fact should share its aggregates if both the conditions below are met:

- Similar non-conflicting dimensionality: If the same fact is used by two virtual data marts, with the same or very close dimensionality, it is preferable to establish

one set of aggregates for both. If however the Product dimension requires “as-is” semantics in one mart and “as-was” in the other, this makes the dimensions conflicting. In this situation if recast-aggregates cannot be avoided they should be separated.

- Access views shall be aggregate agnostic: The ultimate objective behind this requirement is decoupling applications from aggregates. If an aggregate is removed or its level adjusted, applications may need to be tested but should not have to be modified. Access layer views are treated as parts of the application. Their change cascades to the applications that use them. Granted, building access views over aggregate-definition views makes matching and redirection by the optimizer safer. However this also exposes the user or application querying the mart to be conscious of the aggregations. A better design is to establish access views at the base level and capture all idiosyncrasies and computations there. Such views are stable and can be used as the base for all aggregate definition views at higher levels in the lattice. As long as they are straightforward aggregations and all complexity and joins are captured into the base view, the optimizer will detect the subsumption and redirect the queries.

To illustrate the point, Figure 8 shows a base fact with two dimensions combined into a base view. The Base View is used in the two Application Access Views 3 and 4. There are also two system maintained aggregates with their respective view definitions. Access Views 1 and 2 are specific to these aggregates. If applications are exposed to them, then change of aggregate level will require the applications to be changed as well. Decoupling the application entails reliance on access views 3 and 4 only. All queries will be directed to these views, and the optimizer shall redirect to the aggregates whenever it recognizes that the query subsumes the definition of these aggregates. In order to raise the likelihood that this will occur, the base view must absorb all the complexity (filters, computations and joins) and keep the Aggregate Views 1 and 2 as simple as possible. If access views are aggregate agnostic, aggregate levels can be adjusted without requiring applications to be modified, although as part of evaluating the change, their queries must be benchmarked or tested.

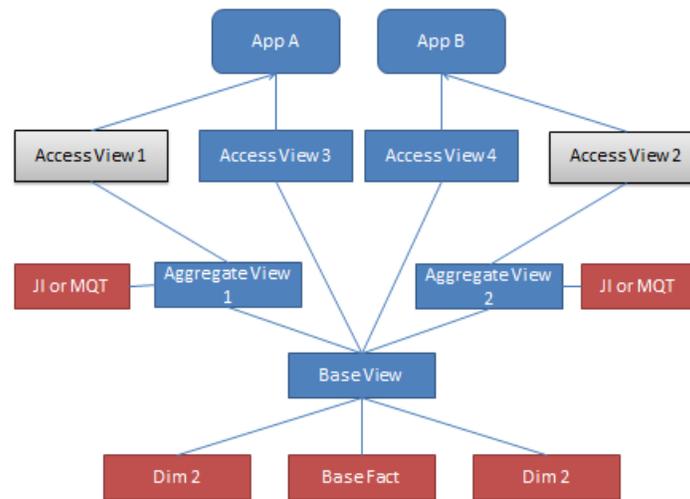


Figure 8 – Application-Aggregate Coupling through aggregate-dependent views

- Materialized aggregates are shared performance objects that belong with the respective base data in the same spirit that indexes are. This idea, however, can only be supported if we decouple applications from those performance enhancers, as is the case with indexes. If we couple performance objects to access layer views we delegate optimal aggregate selection to the application, and any change to these aggregates challenges the “autonomy” of these applications. We consequently propose that aggregate sharing is a virtue if and only if access views and applications are aggregate agnostic.

Automation and Tools

Both during the construction of the lattice and the classification of queries, some level of automation is desirable. Three specific use cases for such automation are:

- Scan actual queries collected from monitoring the system and summarize statistics of tables, views and columns involved.
- Parsing queries to identify their multi-level. This allows us to monitor workloads of queries and know their S-Level, G-Level, M-Level, Q-Level as well as the discretionary and intrinsic cost and rank. This type of analysis is invaluable in analyzing and quantifying performance impact when planning improvement iterations and evaluating results after changes are made.
- Constructing benchmarks from template patterns to evaluate coverage of the lattice, performance improvements, check optimizer choices on use of aggregates etc.

VDM includes customizable tools that can be adapted on a case by case basis. When queries are generated from application tools such as MicroStrategy, the naming

conventions are consistent and simple scanners can achieve a reasonable level of success. To the degree that this level of automation is feasible, and after a potentially tedious initial setup, the repeatability of the process and the opportunity to systematically analyze large samples of actually executed queries makes it worth the effort. Of course the ease of such automation depends on access layer complexity and the consistency of naming conventions

Conclusion

The VDM Aggregation Strategy provides a framework for identifying and refining aggregates and the access layer. From prior applications in other industries, the methodology may yield improvements as high as a level of magnitude. For existing mature systems we typically set a modest target of 20% or 30% improvement of current utilization for targeted periods and workloads. This translates into resource savings that frees up capacity for additional work, and may elongate – or keep from shrinking – capacity upgrade cycles

The process of implementing the VDM Aggregation Strategy is iterative, following these rough steps:

1. Analyze query patterns, set targets and plan
2. Evaluate the access layer design for coupling with aggregates. Evaluate options
3. Create/Refine baseline model envelope; baseline benchmarks
4. Build, prune, load and weigh the aggregation lattice; map current aggregates
5. Analyze historical or projected workloads and generate benchmarks covering the lattice to the extent possible – identify opportunities with existing JIs;
6. Determine optimal aggregates considering existing levels – propose changes and additions
7. Benchmark and evaluate impact
8. Implement
9. Validate

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