

Introduction To Data Mining

(1.1) Motivation & Importance:

The information industry in recent years contains the huge amount of data. This huge amount of data must be transformed into useful information & knowledge. This knowledge & information used in many systems like market analysis, business Management, production & analysis, science & engineering etc., This is the motivation behind the data mining technology.

The evolution of database technology contains the functions like data collection & database creation, data management (store, Retrieve & transaction process) & data analysis & understanding

The database evolution is shown in below diagram.

Data collection & Database creation
 (1960 & early)
 primitive file processing

Data Management Systems
 (1970's - early 1980's)

Hierarchical & network database systems
 Relational database systems
 Data modeling tools: Entity-relationship model etc.
 Indexing & Data organisation techniques:
 B⁺ tree, hashing etc.,
 Query languages: SQL etc.,
 user interfaces: forms, Reports.
 Transaction Management: Recovery, concurrency
 -control, etc.,
 OLTP systems (Online transaction processing system)

Advanced db systems
 (mid 1980's present)
 Advanced data mod-els:
 Object-relational etc;
 Application-oriented;
 spatial, temporal,
 multi-media, scientific

DW & DM
 (late 1980's - pres-ent)
 -DW & OLAP tech-nology.
 -DM & KD

web-base data-base system.
 (1990's present)
 -xml databa-se systems.

Integrated Information systems (IIS)
 (2000 - - -)

fig 1.1 The Evolution of database technology

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In the year of 1960 & early data is stored in the form of primitive file processing. This is used as a base of the rest of the database systems.

In the year of 1970, db systems are developed. The initial db systems are Hierarchical & network, later relational db systems, data modeling techniques, indexing & data organisation & query languages are developed. After that to provide easy understanding to user several user interfaces, form reports & query processing & query optimization, & transaction management techniques are developed.

In the year of early 1980's OLTP (Online transaction processing systems are developed).

In the year of mid 1980's advanced db systems are developed like advanced data model, application oriented db systems are developed.

In the year of late 1980's Data warehousing DW has been developed. It is a repository of heterogeneous data sources organised and

etc.

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a schema to support management decision.

Data warehouse contains the OLAP (Online analytical processing) technology. OLAP contains the functions like summarization & aggregation. Later on data mining & knowledge discovery techniques are developed.

In the year of 1990 web based db systems are developed. By integrating above three systems, we get integrated information systems. This systems developed in the year of 2000.

Therefore, the information industry contains the large amount of data & also data is stored in different sources that exceeds human ability. Therefore, we need data mining tools to analyse & search for interesting patterns or knowledge. This knowledge is used by the management to take accurate decisions.

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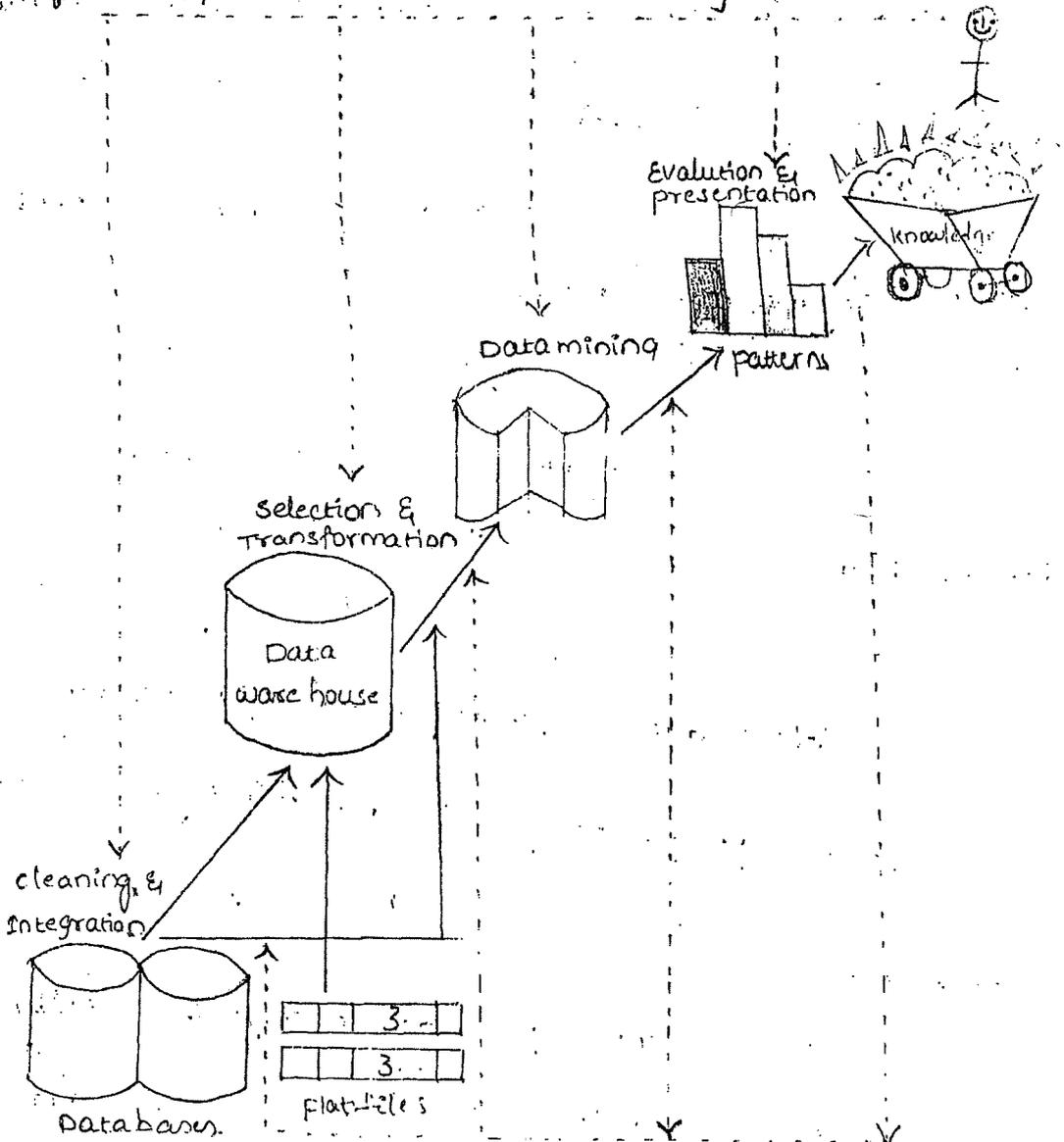
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1.2 what is data mining:

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Data mining refers to extracting or mining data from large databases. Some of the people think that data mining is a synonym of KDD knowledge discovery in databases & also some of the people think that data mining as a essential step in the the process of KDD. This is shown in below diagram.



Data mining as a step in the process of knowledge discovery

1. Data cleaning:

The data from the large heterogeneous databases are retrieved & noisy data or inconsistent data is removed before going to the next step.

2. Data Integration:

The data from the different sources are integrated & then this data is loaded into data warehouse

3. Data Warehouse: (DWH)

It is a centralized repository it contains all the organizations data.

4. Data Selection:

The selected data is retrieved from the data warehouse.

5. Data Transformation:

Here data is transformed into a form that must be appropriate for the data mining.

6. Data Mining:

It is essential process. It contains intelligence -nt techniques or methods like summarization, aggregation etc.

7. pattern evolution:

databases

Here interesting patterns are evolved.

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8. knowledge presentation:

Here visualization techniques are applied

to present mined knowledge to the user.

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(Based on this the typical data mining

warehouse

architecture contains the following components.

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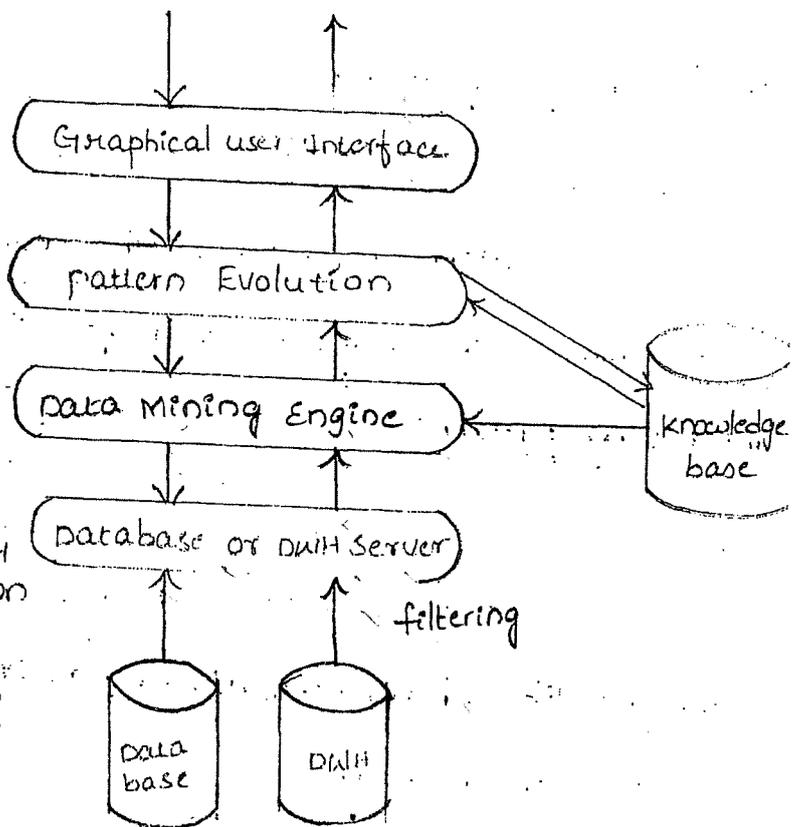


fig 1.3 : Typical data mining Architecture

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1. Database or DWH or Any other information Repository:

Here data is retrieved from any data base, or data warehouse or any other information repository we apply data cleaning & integration technique before data is going to database or DWH server.

2. Database or DWH server:

It contains all the data according to the user specification.

3. Knowledge base:

Here we retrieve the domain knowledge & this domain knowledge is used for searching the interesting patterns.

4. Data Mining Engine:

It is an essential process. It contains technique like characterization, classification, association & cluster analysis etc.

5. Pattern Evolution:

Here interesting patterns are evaluated.

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6. Graphical user interface:

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It provides the communication between

user & data mining system through this end user can

present the queries to data mining system.

1.3 Data Mining - on what kind of data?

Data mining can be applicable to any

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information repository. This information repository may

be a relational databases, DWIT, transactional database

advanced database systems, & Advanced database appli

-cations.

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* 1.3.1 Relational Databases:

Relation database system consists of tables

& also each table contains the unique name. Each

table consists of set of attributes or columns &

generally stores large set of data in the form of

tuples, or records or rows. As well as each row is

uniquely accessed by using a key.

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association

Example:

3.

For example, relation database tables for ALL Electron-

ics Company contains the tables like customer, item,

employee, purchases etc.,

Customer

cust-id	name	address	age	income	...
C1	Smith	123 Ham-st Canada	21	\$25,000	

Here customer table contains the attributes like cust-id, name, address, age, income etc.,

Item

Item-Id	name	brand	category	type	place-made	cost
I3	high res-TV	Toshiba	high resolution	TV	Japan	\$2000

Employee

emp-id	name	category	group	salary
E5	John	home, entertainment	manager	\$10,000

Branch

Branch-id	name	address
b1	city square	123 main st, Toronto, Canada

purchases

trans-id	cust-id	item-id	date	time	cost
T100	C1	I3	29/06/09	15:45	\$100

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Item_Sold

Trans_id	Item_id	quantity
T100	I3	1

fig 1.4 Relational tables for AllElectronics database

Here some of the tables represents the relation between multiple tables like purchases, item-sold etc.,

The Relation database is a most popularly available & rich information repository in data mining. In Relation databases data modelling is done by using ER-model

1.3.2 Data Warehouse (DWH):

It is a centralized repository organised under a schema to support management decision. We load data into 'data warehouse' i.e., the data warehouse is constructed by data cleaning, data transformation, data integration & data load.

for example, typical data warehouse for All Electronics company.

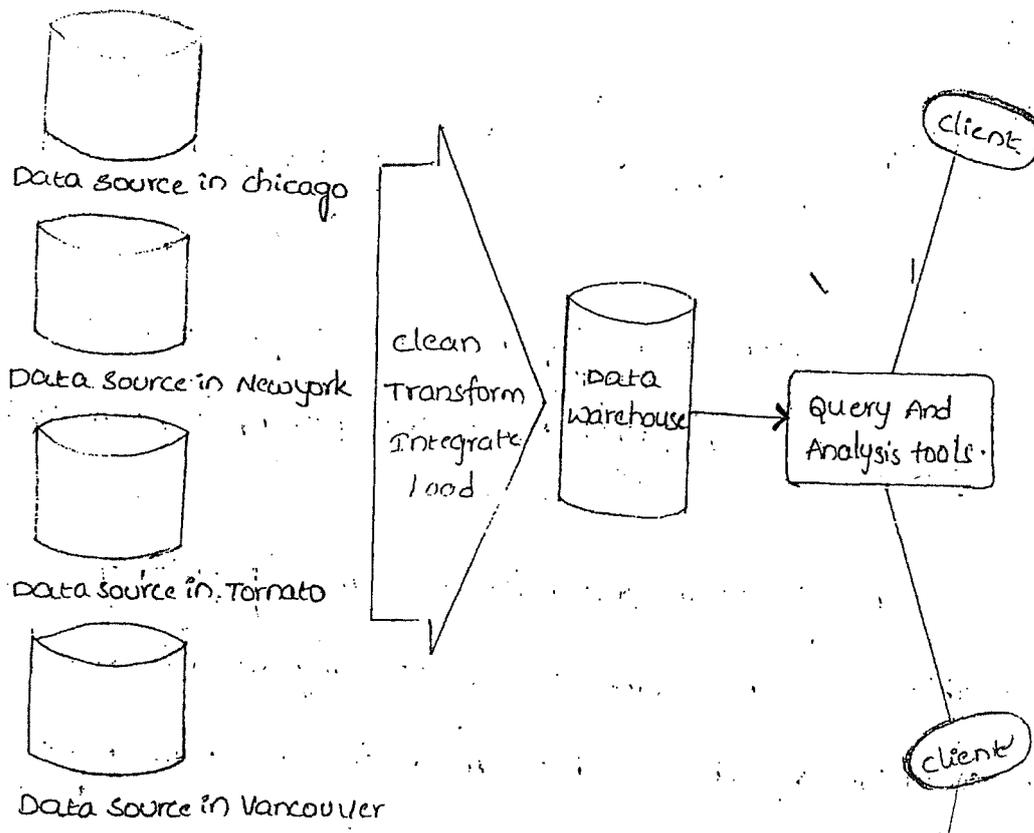


Fig 1.5 Typical Architecture of Data Warehouse for All Electronics company.

Here All Electronics company has branches All around the world. Therefore data from different branches first of all cleaned, transformed, integrated & finally loaded into data warehouse, then we apply query & analysis tools before going to present the data to the management.

Data warehouse contains the historical data

e.g. 5 to 10 years of data & used to support

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management decision. In data warehouse data modeling is done by using schema but the physical architecture of data warehouse contains the "data cube". Through this data cube "multi-dimensional view of data is presented."

Example

Data cube for summarized sales data of "ALLElectronics" is shown below. This data cube contains the three dimensions i.e., address (it contains city values Chicago, New York, Toronto, Vancouver), Time (Quarter values Q1, Q2, Q3 & Q4) & item (it contains item type like home entertainment, computer, phones, security).

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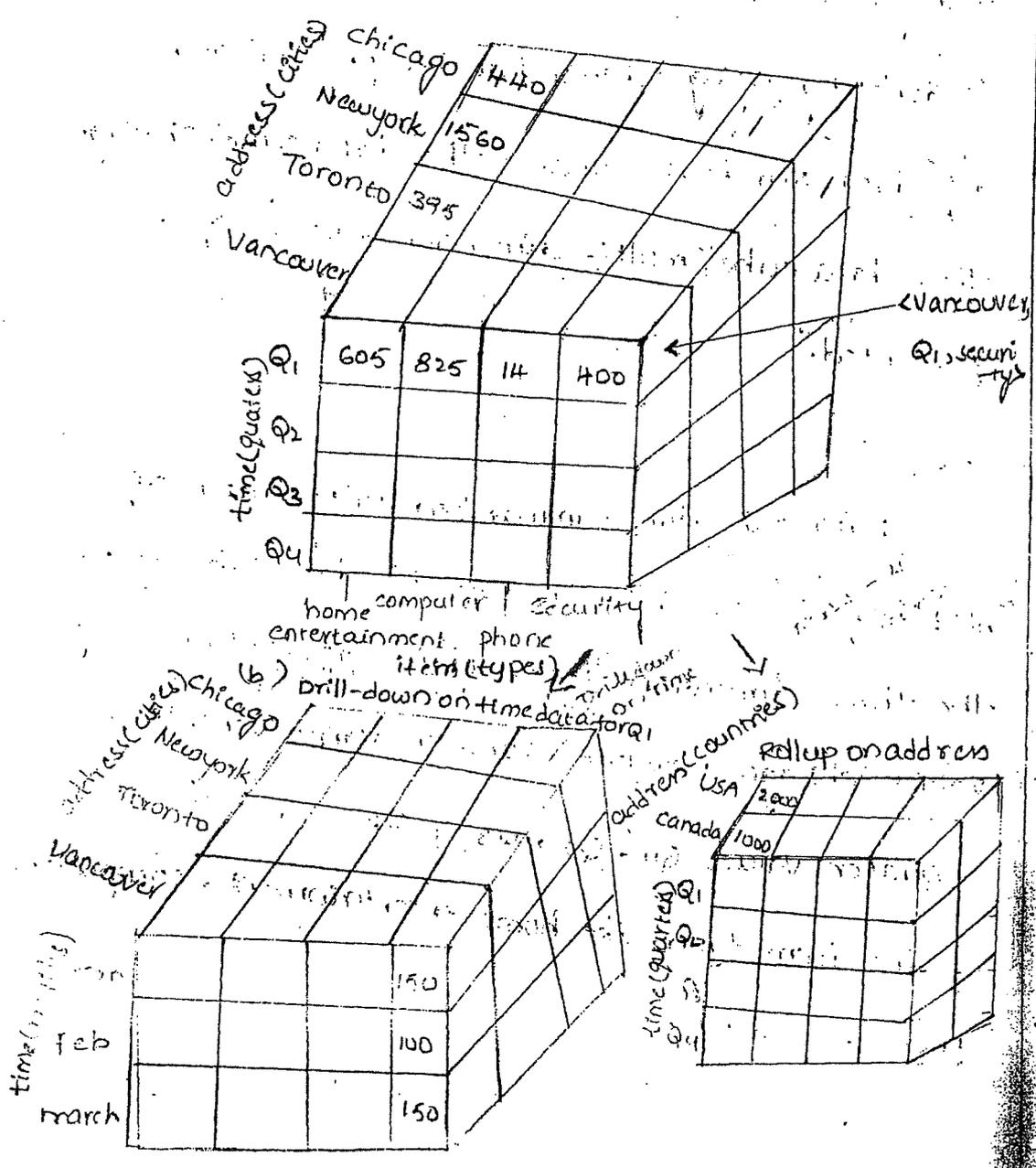


fig 1.6 multi-dimensional data cube for AllElectronics

- 1.6 (a) : summarized sales data
- (b) : Drill down & roll up operations

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Here data warehouse contains the data about all the subjects & all the organizations. Therefore, its scope is enterprise-wide. On the other hand datamart, it is a subject of data warehouse. It contains the single subject area & its scope is department-wide.

Data warehouse provides multi-dimensional view & summarized data. Therefore, data warehouse is well suited to OLAP (Online Analytical processing).

OLAP contains the techniques like

"drill-down" & "roll-up"

In fig 1.6 (b) drill-down on the time data for Q₁ by monthly & also rollup on address by countries.

Electronics

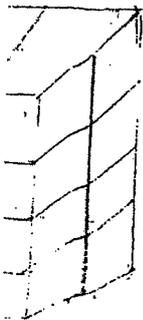
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Transactional Databases:

Transactional databases contains the files where each record represents a transaction. A transaction uniquely identified by the transaction-id

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address



trans-id. These transactions are can be stored in tables & each record/row represents transaction.

The fragment of transactional database for

All Electronics is shown in below.

Trans-id	list of item-id
T100	I1, I3, I8
.....

Fig 1.7 Fragment of Transactional database for sales at All Electronics.

Here the sales table contains the nested relation i.e., it contains the list of item-id's

1.3.4. Advanced Database systems & Advanced database applications:

The new applications like spacial data (such as maps), engineering & design (such as designing buildings, designing components, Integrating circuits), Text & multimedia (such as audio, video, images) Temporal & time related data (such as historical data & stock exchange data) & world wide web (widely distributed information repositories on the internet) etc,

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Object-oriented database systems:

These object oriented database systems are mainly based on object oriented paradigm. Here each entity is treated as object, in this the data & the code related to object are encapsulated into single unit.

Each object contains the following.

(i) set of variables that describes the object:

This is similar to attributes in ER-model

(ii) set of messages to communicate with one object to another object.

(iii) set of methods each method contains the code to implement message.

The similar objects are combined into class.

This class is known as object class then each object of this class is an instance for that class.

Eg: employee class contains variables like name, address, salary etc.

Object-related database systems:

The object related database systems are constructed from object relational data model.

These models are extended from relational model & also new data types has been added to handle the complex objects. The constructors also defined to handle this added data types in relational query languages.

In data mining system the object oriented database systems & object related database systems share some similarities.

spacial database systems:

spacial database systems contains the data in the form of geographic maps & are represented in the form of Raster format with n-dimensional bit maps or pixel maps.

For example, 2D satellite image is represented in raster format, where each pixel represents the rain fall in this specific area.

The maps are also represented in vector form i.e., where buildings, roads etc, are represented in the form of points, polyline or polygon.

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Text & multimedia Database Systems:

Text database systems contains long sentences, paragraphs to specify product specification or error or bug specification or summary report specification etc., These text database systems are unstructured.

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Multimedia databases contains the data in the form of audio, video, images. These are used in applications like voice message systems, video on demand systems, world wide web information systems etc.,

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Temporal & Time-series Database systems:

These two database systems contains the set of data. This data continuously changes with time like stock exchange data.

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World Wide Web information systems:

These are the widely distributed information systems like yahoo or America online or online service. These are linked with objects to exchange the information through internet i.e., the end-

user can get these services through internet & also end user get attractive web pages. These web pages are un-structured i.e. doesn't contain any schema or pattern.

★ List Data Mining functionalities:

Data mining functionalities are used to specify the kind of patterns to be mined found in data mining tasks.

Data mining tasks mainly classified into

1. Descriptive

2. Predictive

"Descriptive means it explains the general characteristics of a data in database"

"Predictive means it provides conclusion on current data through prediction"

Here end user can go for the search for multiple patterns because end user does not know

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(What kind of patterns are interesting.)

Therefore, data mining system provides the

multiple search facility & also provides the interesting patterns are search in various ^(levels) granularities.

Therefore, data mining system provides the

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hints to the end user to search for the interesting patterns.

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The data mining functionalities mainly classified

1.4.1 concept / class description: characterization & discrimination:

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The data is associated with the concept or a class. for example in ALLElectronics database class of items for sale contains the computers & printers.

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similarly concept of customers include big spenders & budget spenders.

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Therefore, the data associated with concept or class is called as concept / class description. This

descriptions are derived via

1. Data characterization
2. Data discrimination
3. Both data characterization & data discrimination.

1. Data characterization:

It is the summarization of general characteristics or features of a target class.

Data characterization can be achieved through several techniques.

for ex, Rollup operation in OLAP.

once the data characterization is completed then data can be presented in various forms "bar charts, pie charts, multi-dimensional data cubes"

2. Data Discrimination:

It is the comparison of target class with one or more constructing classes. This is called as the "data discrimination".

Here End user specifies the target classes & constructing classes & also End user can

the data from database by using queries:

3. Data characterization & Data Discrimination:

Note: we have to write both (1) & (2) here.

1.4.2 Association Analysis:

The association analysis is can be used to identify the association rule between each pair of attributes.

This association analysis mainly used in transactional data analysis.

The association rule $x \Rightarrow y$ i.e.,

" $A_1 \wedge A_2 \dots \wedge A_m \rightarrow B_1 \wedge B_2 \dots \wedge B_n$ " where

A_i for $i \in \{1, \dots, m\}$ and B_j for $j \in \{1, \dots, n\}$ for pair of attributes.

For example, All Electronics database contains the association rule

$\text{age}(x, "20 \dots 29") \wedge \text{income}(x, "20k \dots 30k") \Rightarrow \text{buys}(x, "cd player")$

[support = 2%, confidence = 60%]

as mentioned in the above ex, a person whose age between 20 to 29 & whose income will be 20k

to 30k can buy or purchase a CD player at ALL Electronics Company. With support 2% & confidence 60%.

(Or)

This association rule indicates that $x \Rightarrow y$ can be interpreted as customers 2% support are 20 to 29 years of age with income 20k to 30k buys or purchases a CD player at ALL Electronics Company.

There is a probability confidence 60% of this age group & this income can buy CD player.

1.4.3 classification & prediction:

The classification is the process of finding set of models or functions that describes class i.e., the classification mainly used to identify class labels. But many of the applications the end user search for the data. In most of the cases usually we search numerical data.

Therefore search for numerical data

fig 1.8

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in database is called "prediction".

1.4.4 Cluster Analysis:

Cluster Analysis is nothing but identifying the data objects with respect to our requirement; we have to also examine the data object which are not relevant to our class label.

By performing cluster analysis, we can easily identify similarity & dissimilarity of class label. So that we can group the similar objects & we can leave the non-similar objects which is represented in below diagram.

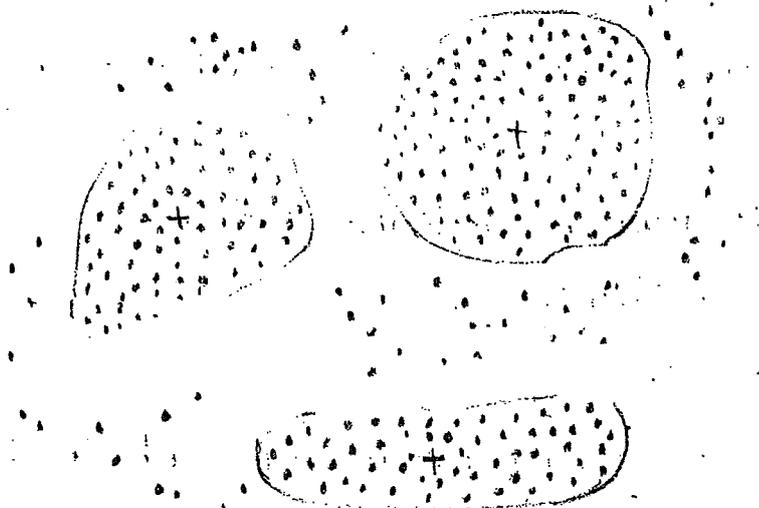


fig 1.8

A 2-D plot of customer data with respect to customer locations in a city, showing three data clusters. Each cluster "center" is marked with a "+"

1.4.5 Evolution Analysis:

The data evolution analysis describes data objects that are changing with time.

These data objects are evolved through characterization, discrimination, association analysis & cluster analysis.

1.5 Are All of the patterns interesting? :

(Or)

Interestingness of a pattern:

The datamining system capable of providing 1000's of patterns. A pattern is interesting if

* That must be understandable to the human beings.

* It must be applicable for new data.

* potentially useful.

* Novel.

Interesting pattern represents the knowledge. Several objective measures for interestingness of a pattern. In these association rule is one of the

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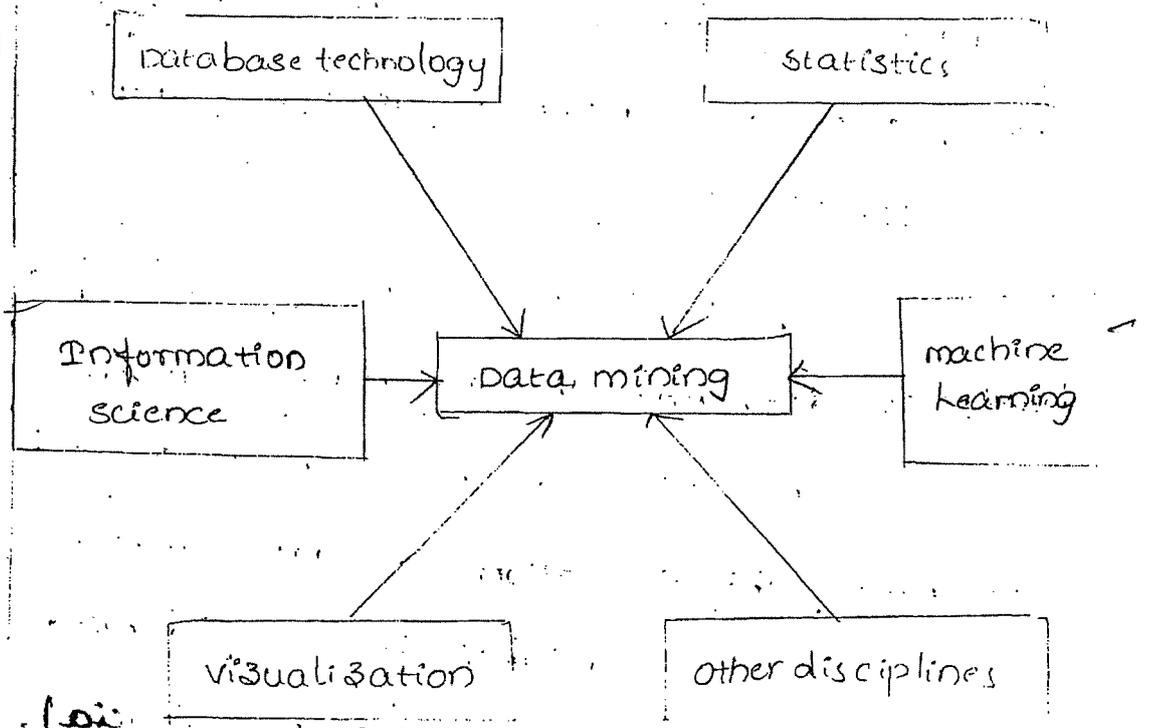
method. Therefore association rule $(x \Rightarrow y)$ then support $(x \Rightarrow y) = P(x \cup y)$, where $x \cup y$ indicates that a transaction contains both x & y i.e., the union of item x & item y .
 Confidence $(x \Rightarrow y) = P(y/x)$ conditional probability i.e., the probability that a transaction containing x also contains y .

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1.6 Classification of data mining systems:

The classification of data mining system is an integration of multiple disciplines (fields). This is shown in below

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fig 1.9 classification data mining systems

Here data mining system is integrated with database Technology, statistics, machine learning, visualization, information science etc.,

Therefore data mining system is classified as following

1. Classification According to kind of database it is mined:

The data mining system is classified based on database mined. But the database technology is classified based on data model or type of data (i.e., application oriented).

If the database technology is classified based on data model this may be a relational database, transactional database, object-oriented database, object relational or data warehouse.

If the database is classified based on the type of data then it may be a spacial, temporal & time series, text & multi-media etc.,

Systems.

2. classification according to the kind of knowledge it is mined:

Data mining system is classified based on knowledge mined i.e., we use the data mining functionalities like concept or class description, association analysis, cluster analysis & evolution analysis.

Data mining system is also classified based on knowledge mined at different levels. or top level or primitive level (low level or raw level) or multiple levels.

3. classification according to kind of techniques utilized:

The data mining system is classified based on techniques utilized. These techniques describes the degree of user interaction involved.

For example query-driven systems, autonomous systems, interactive systems etc., or methods for data analysis.

Classification according to kinds of applications

adapted:

The datamining system is classified based on applications adapted like finance data base system, tele communication database system, share market database system etc.

1.7 Major issues of data mining:

These are mainly classified into

1. mining methodology & user interaction issues.

2. performance issues.

3. Issues related to database types.

1. mining methodology & user interaction issues:

These specifies the kinds of knowledge mined, knowledge mined at different levels,

use of domain knowledge & knowledge presentation.

Mining knowledge from databases:

different users require different

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kinds of knowledge. Therefore, data mining system must provide wide range of data analysis & also knowledge discovery through different data mining techniques, like data characterization, data discrimination, association analysis, cluster analysis & evolution analysis.

1.2 mining knowledge at different levels:

issues:

Here end user interact with the data mining system & use the different ^{Online analytical processing} OLAP techniques like drill down & roll up, through these techniques end user mined the knowledge at different levels.

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1.3 Background knowledge:

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This background knowledge guides the discovery process if end user has a background knowledge about the database like the constraints or association rules or conditions then this data mining process is speed up.

1.5 presentation & visualization

The mined knowledge is presented to the end user & this must be understandable to the end user. Therefore, we use the several visualization techniques like trees, graphs, charts, matrix etc.

1.4 Data mining query languages for data mining

Tasks:

like the relational query language, for example SQL. In SQL we present ^(temporary query) Adhoc Query & get the data. Similarly data mining query languages has to be developed & it must support end user Adhoc data mining tasks, like data analysis, get the domain knowledge, understand the constraints & conditions etc.

1.6 Handling noisy (or) incomplete data

Generally databases contains the noisy or incomplete data. Therefore, we use the data mining techniques like data cleaning &

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data analysis to avoid this noisy or incomplete data.

1.7 Pattern Evolution:

still the data mining system uncover thousands of patterns & also many of the discovered patterns uninteresting to the user. Therefore, end user has to specify the measures or constraints to reduce the search criteria.

2. Performance Issues:

These issues contains the efficiency, scalability & parallelization of data mining algorithms.

2.1 Efficiency & scalability of data mining

Algorithms:

We have to retrieve the data from large databases. Therefore, the data mining algorithms must be efficient & scalable.

In data mining system knowledge discovery, efficiency & scalability are the main

key terms.

2.2 parallel, distributed & Incremental data mining

algorithms:

End user has to access the data parallelly from different data sources & also data is distributed to different data sources. These two are done by using parallel & distributed data mining algorithms. In these algorithms data is divided into partitions then partition processed parallelly & then results of this partitions are merged.

The incremental data mining algorithms are mainly used to reduce the cost of data mining process & also update the data in database without performing the search from the beginning of Database.

3. Issues related to database types:

3.1 Handling relational & complex data:

The relational data is handled

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efficiently by using relational database systems & data warehouse systems.

parallelly

The complex data like spatial data, text

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& multi-media data, time series data. These

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complex data is handled by using systems spacial

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3.2 Mining data from heterogeneous databases

Mining the data like LAN systems,

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WAN systems, distributed systems & Heterogeneous database systems is only possible through data mining

of data

System. This data mining system also provides &

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improves the information exchange & interoperability

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in heterogeneous databases.

Chapter - 2

Data Warehouse And OLAP Technology for Data Mining

Basic concepts of data warehouse:-

(i) What is Data Warehouse

According to warehouse Inman, "A Data Warehouse is a Subject oriented, integrated, non-volatile, Time-Variant collection of data, to support management decisions".

Subject Oriented:

For example, consider the retailer, these retailer contains the Database systems like Retail Database System, Catalog Database system and Outlet Database Systems. These database systems individually support for different queries. But a user want to run a query in all the sales. This is only possible through data

- warehouse. therefore Data warehouse organizes Subject areas like customer, products, suppliers, Sales etc. Here the subject area is "Sales". This is shown in the following diagram.

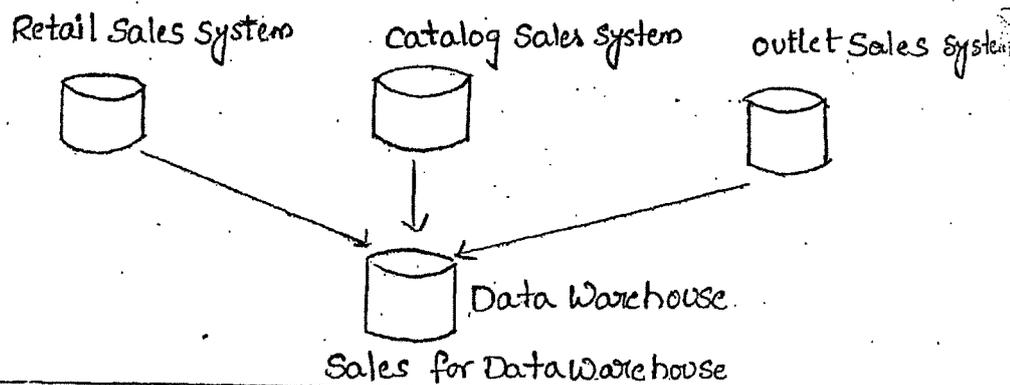


Fig C2.1.1) Subj-Oriented Sales information

Integrated :

Here, different data sources data are integrated. Then this data loaded it into Data Warehouse. These sources may be relational databases, flat files, Excel sheets, Online Transaction records, etc. Before the data is loaded it into DWH, we apply data cleaning and data analysis.

In the above example, 3 data source, data are integrated then we get the unique key. Using this unique key, we uniquely identify the data record in datawarehouse. This is shown in the following diagram.

Sys

Retail Sales System

Catalog Sales System

Outlet Sales System

Product-code:

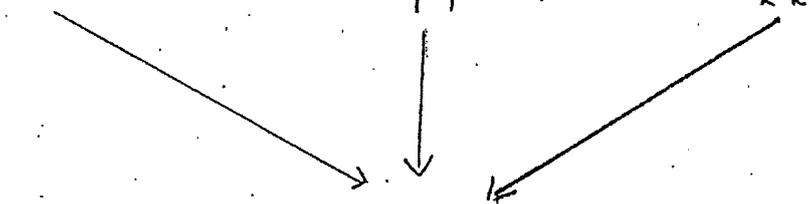
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Product-code:

yy

Product-code:

zz



Product-code for DWH:

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Fig (2.1.2) Integrated information for Sales

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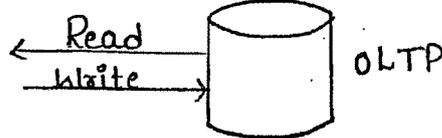
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Non-Volatile:

The non-volatile means read only. The DWH User always read the data. But the OLTP user can read the data as well as write the data. This is shown in below.

USER



USER

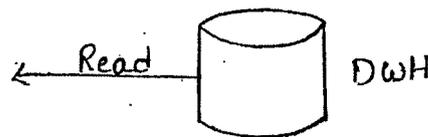


Fig (2.1.3) Non-Volatile

Time - Variant:

The DWH contains the historical data.
i.e; 5 to 10 years of data. But the OLTP System
contains the current data.

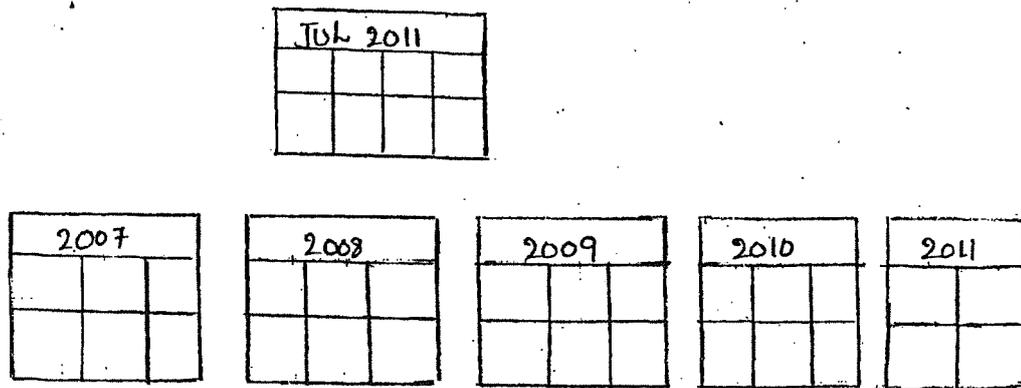


Fig (2.1.4). Time Variant

Finally we extract the data from the DWH & then we apply the different visualization Techniques like Trees, Graphs, charts, Tables, etc. to present the data to management. The management using this data they take the accurate decision.

2.1.1 Differences between operational database systems & Data warehouses:-

The online operational database systems mainly used for Online transactions & Query processing. These include the transactions like sales, payroll, marketing, manufacturing, etc. These type of systems are called as "online Transaction Processing" (OLTP) Systems.

These systems mainly used for day to day transactions.

The DWH contains the historical data. This is used for data modelling and data analysis. This is used for decision making. This type of systems are called as "online Analytical Processing" (OLAP) systems.

Advantages of Data Warehouse:-

1. It makes the data permanent.
2. It makes the data accessible.
3. It identifies hidden business operations.
4. It improves the customer relationship.
5. It provides the security.

Differences between OLTP and OLAP:-

1. Users and System Orientations:

The OLTP system is customer orientation to support the Online transaction.

The OLAP system is market orientation to support decision making.

2. Data: The OLTP systems contains the current data. This data does not be used for decision making.

The OLAP system contains the historical data and supports functions like summarization & aggregation and also data stores and manages at different levels.

3. Database design: In OLTP systems data modelling is done by using ER-model.

In OLAP systems data modelling is done by Star (3) Snowflake scheme.

4. View: Using OLTP Systems we extract the data within a Enterprise & department.

In OLAP systems we access the data from different organization's & different data sources.

5. Access pattern: The OLTP systems contains the data. This data used for online transaction's..

The OLAP systems contains the historical data. This data is used for data Analysis.

DWH → OLAP
OpDBS → OLTP.

Users and system orientation
Data
Database design
view
Access pattern.

Comparison between OLTP and OLAP:

Feature	OLTP	OLAP
Characteristic	Transactional - Processing	Informational - Processing.
Users	Clerks, OBA, DB - Professional	HR, manager, Analyst, executive.
Functions	Day-to-day transa- -ctions.	Decision making.
Orientation	Online Transaction - orientation	Analysis Orientation
Data	Current data	Historical data
Database Design View	ER- model Relational (2-D)	Star (3 rd) Snowflake Schema Multi Dimensional.
Units of work	Simple Transactions	Complex Queries.
Access Pattern	Read / Write	Read only
Number of Users	Thousands	Hundreds
Number of data - records accessed	Tens	millions
Data storage	100MB to GB	100 GB - TB

Fig (2.15). Comparison between OLTP & OLAP

Data Warehouse Modeling: Data cube & OLAP

2.2. Multi-Dimensional Data model:

The Data Warehouse (or) OLAP tools contains the data in the form of multi-dimensional model. i.e; It contains the Data cube.

2.2.1. From Tables and Excel sheets to Data cubes:

Using Data cube we can view the data (or) analyze the data in the form of multi-dimensional model. It is defined by Dimensions and Facts.

The Dimension is nothing but the Entity. Using this organizations store the data. For example, sales DWH contains the dimensions like Time, Item, location, supplier etc. Using this dimensions we can analyze the things like monthly sales of a item, branches & locations at which the item were sold.

This Dimension information is stored in a Table. This is called as "Dimension Table". Each Dimension Table contains the Set of Attributes. For example, Item-Dimension contains the attributes

like item-id, name, category, brand, type, etc.

The multi-dimensional data model the entire data typically organized under the central theme like sales. This "Central Theme" is called "Fact Table".

The facts are numerical entities (or) numerical measures like dollars - sold (Sales amount in dollars), Units - sold (Total units sold).

For example, consider the table for sales DWH for All electronic's company is shown in below. It contains the dimension's like Time, items, and location.

Location = "VanCouver"				
Time (Quarter)	Item (types)			
	Home Entertainment	Computers	phones	Security
Q1	600	700	500	400
Q2	700	800	600	700
Q3	500	600	400	700
Q4	600	700	800	900

Fig (2-2-1-1). Table for sales DWH, it contains Dimensions, time, item & location

The table information is represented in 3D-cube
 The 3D-cube is shown in below.

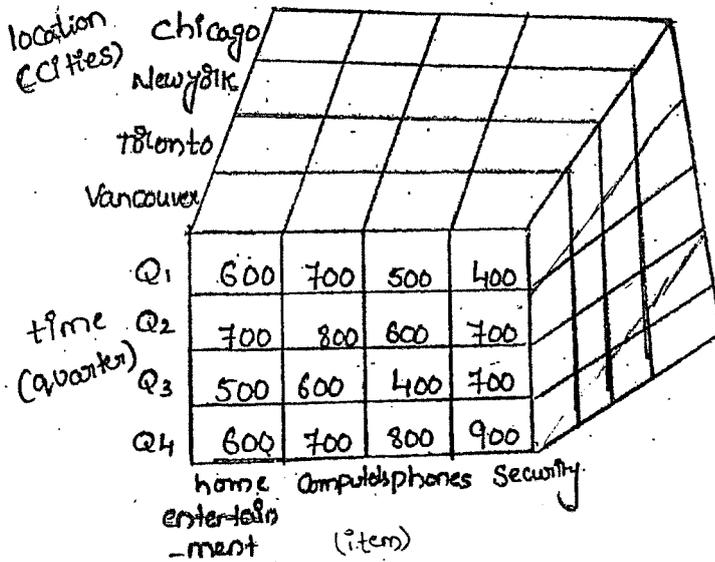


fig (2-2-1-2): A 3-D data cube for sales
 it contains dimensions time, items & locations.

The Datacube contains the n-dimensions.
 The above Data cube may contains the set of dimensions. Using this set of Dimensions we can construct the ^{relation} lattice of cuboids. Each cuboid gives the different level of summarization. Once the cuboids are completed we make the data cube. This is shown in the following diagram.

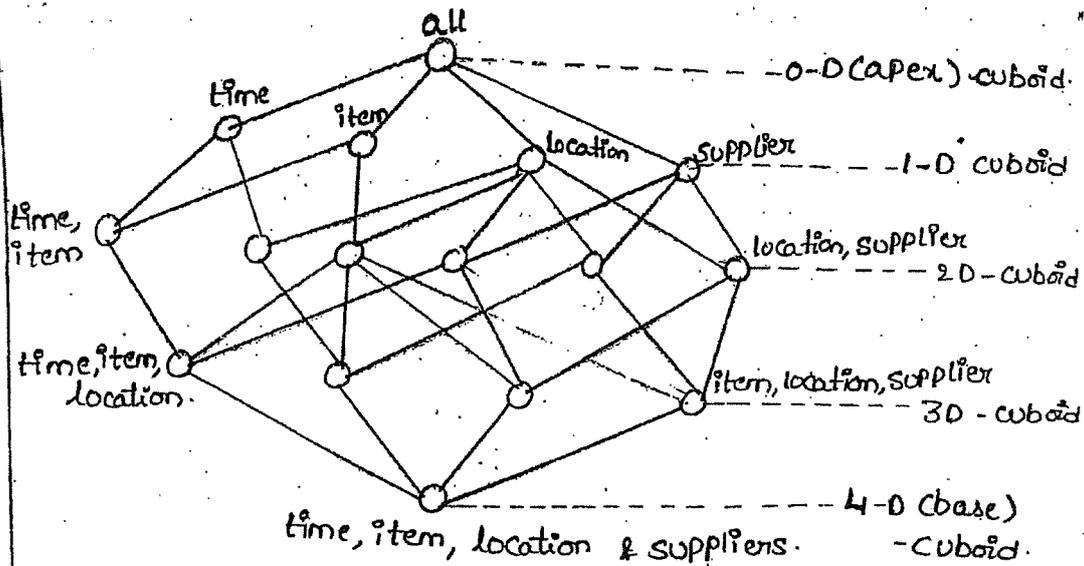


Fig (2.2.1.3): A lattice of cuboids makes 4D Data cubes. It contains Dimensions time, item, location & suppliers.

In the above cuboid, 4D cuboid is called as the "Base cuboid". It gives the lowest level of summarization. The 0-D cuboid is called as "apex" cuboid. It gives the highest level of summarization. It is represented by the keyword "all".

2.2.2. Star, Snowflake and fact constellations

Schemas for multidimensional databases:-

In Relational database the data is loaded by using entity Relationship model. This is best suitable for the online transactions. But the Data Warehouse contains the subject areas like customers, products, sales, suppliers, etc. This is used for data analysis.

Therefore to load the data into DWH we use multi-dimensional Data model. Such model exists in form of star schema (or) snowflake schema (or) fact constellation schema.

Star Schema:

That is the most popular data model to load the data into DWH. It contains 2 tables.

1. Fact Table (It contains large amount of data without any duplication).
2. Dimension Tables (one table for each dimension)

This Schema look like the star. Because of that it is named as "star schema". This is shown in below.

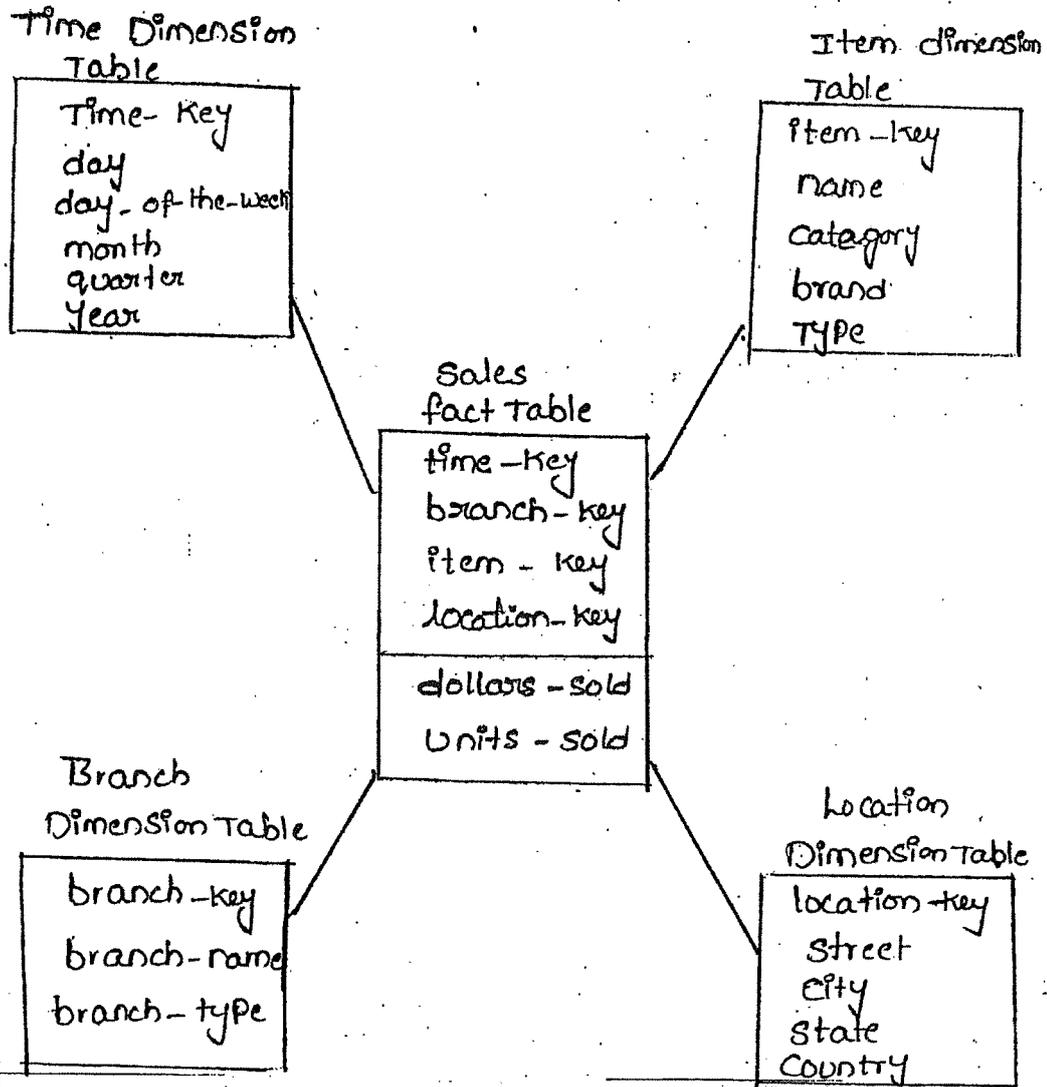


fig (2-2-2-1): Star Schema for All Electronics

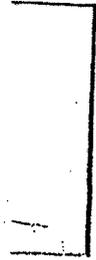
Sales DWH

Star schema for fact Table contains 2 parts.

- 1) Key for each dimension table.
- 2) Measures that is to be analyzed.
i.e; like dollars-sold, units-sold.

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Dimension



Table

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Snowflake Schema:

In Star schema the dimensions tables are not normalized. But in snowflake schema the dimension tables are normalized.

i.e; The tables are further splitted it into more tables. This is shown in below.

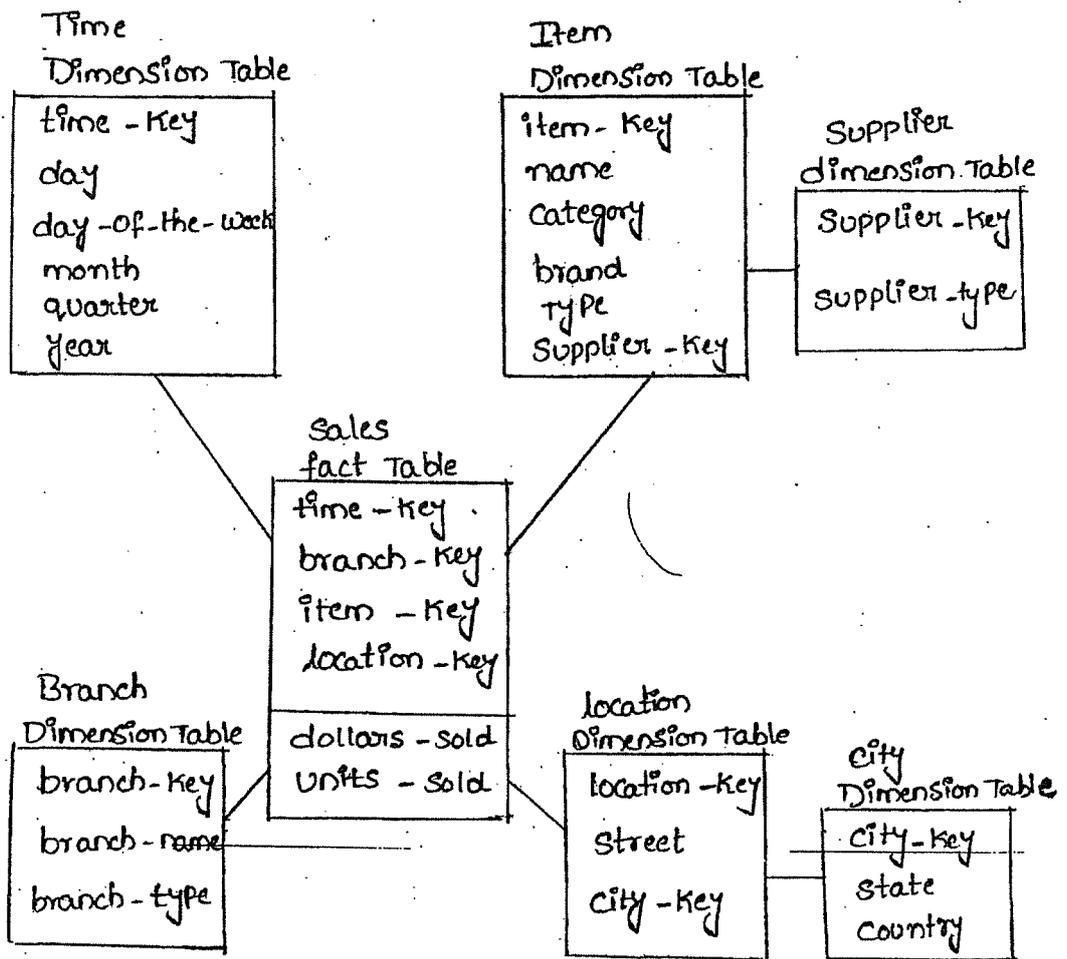


Fig (2.2.2.2) Snowflake Schema for AllElectronics Sales DWH

Fact Constellation Schema:

In some of the applications it requires the multiple fact tables to represent the Dimension Tables. Such a schema is called as the "fact Constellation Schema".

The DWH contains the information about the entire the organization. Its scope is enterprise wide. But the Data mart is subset of DWH and it contains the single subject area its scope is department wide. If it is a DWH (or) Data Mart - it contains 2 Definition's.

(i) Cube Definition:

Syn: $\text{define cube } \langle \text{cube-name} \rangle [\langle \text{dimension-list} \rangle]$
 $\langle \text{measures-list} \rangle$

(ii) Dimension Definition:

Syn: $\text{define dimension } \langle \text{dimension-name} \rangle \text{ as } \langle \text{attributes-list} \rangle$

2.2.3. Measures:-

These are the numeric values, that is to be analyzed. we find the measure value at anytime

by aggregating the data with corresponding dimensions.

The measures are mainly classified into 3.

(1) Distributive:

If it is a distributive measure, then it contains the function's like $\text{sum}()$, $\text{max}()$, $\text{min}()$, $\text{count}()$, etc....

(2) Algebraic: If the measure is algebraic means it contains the function's like $\text{avg}()$. The average is

$$\text{avg}() = \frac{\text{sum}()}{\text{count}()}$$

Here $\text{sum}()$ & $\text{count}()$ are distributive measures.

(3) Holistic:

If the measure is holistic means it contains the functions like $\text{mean}()$, $\text{median}()$, $\text{mode}()$.

2.2.4. Introduction to Concept Hierarchies:-

The concept hierarchies defines sequence of mappings from lowest level to highest level. The concept hierarchy for the location is shown in the following diagram.

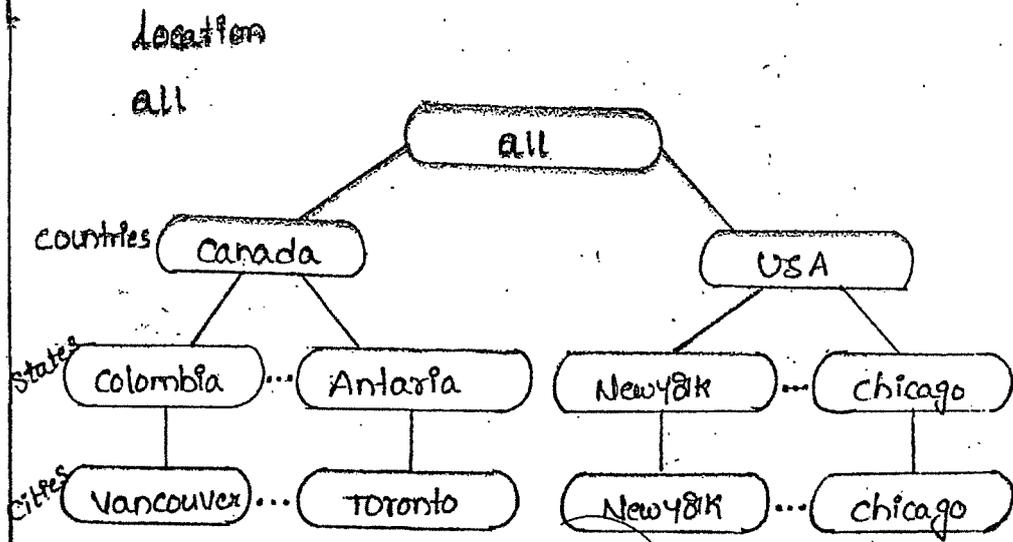


Fig (2.2.4.1): Concept hierarchies for location

Dimension

The concept of hierarchy for location dimension contains the attributes street, city, state and country. Using this attributes we can define the concept hierarchy.

" Street < city < State < country "

These attributes are organized in partial order then we get the "lattice". The lattice for location

Dimension.

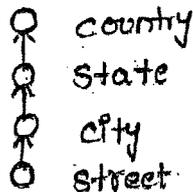


Fig (2.2.4.2). Lattice for location Dimension

We can also define the lattice for time Dimension.

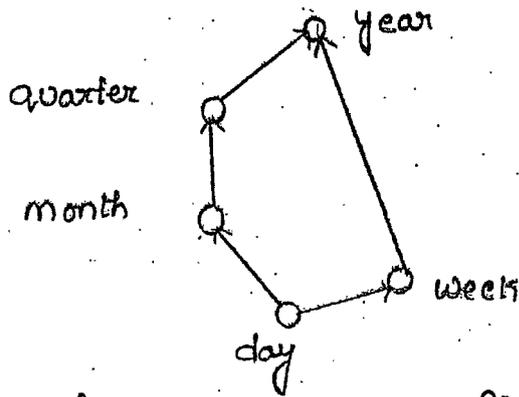


fig (2.2.4-3). Lattice for Time Dimension

The concept hierarchy is also defined for grouping of values then it is called as "Set-grouping hierarchy". The Set Grouping hierarchy for price Dimension is shown in below.

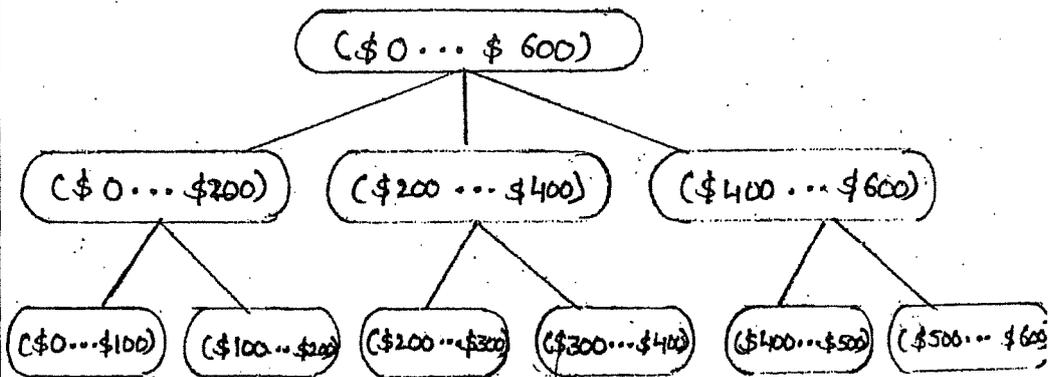


fig (2.2.4-4). Set Grouping hierarchy for price dimension

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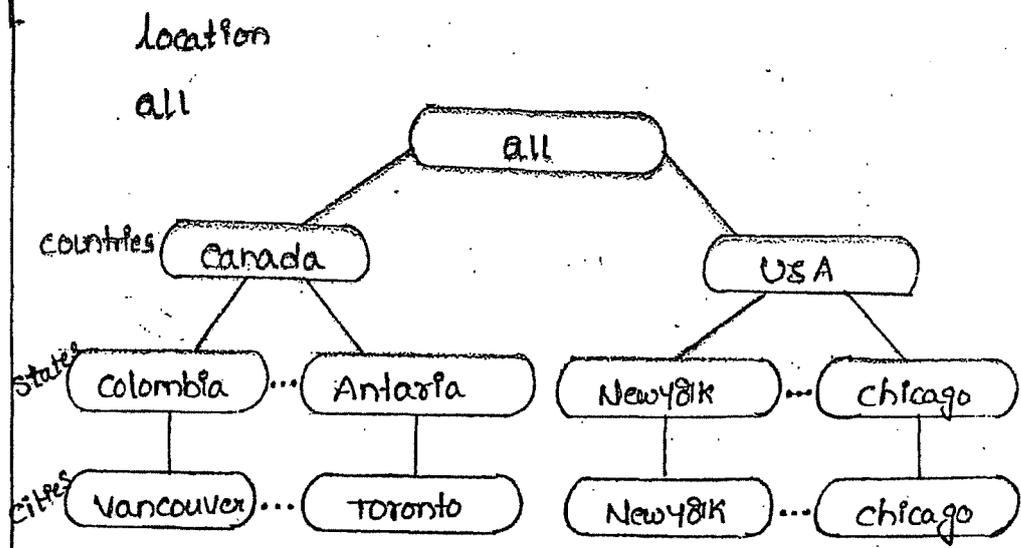


Fig (2.2.4.1): Concept hierarchies for location
Dimension

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"Street < city < state < country"

These attributes are organized in partial order then we get the "lattice". The lattice for location Dimension.

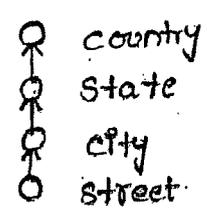


Fig (2.2.4.2). Lattice for location Dimension

We can also define the lattice for time Dimension:

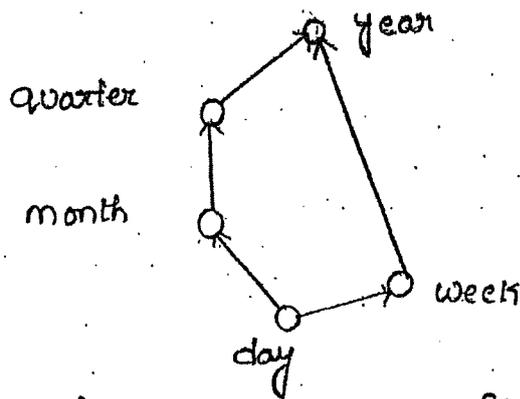


fig (2-2-4-3). Lattice for Time Dimension

The concept hierarchy is also defined for grouping of values then it is called as "Set - grouping hierarchy". The Set Grouping hierarchy for price Dimension is shown in below.

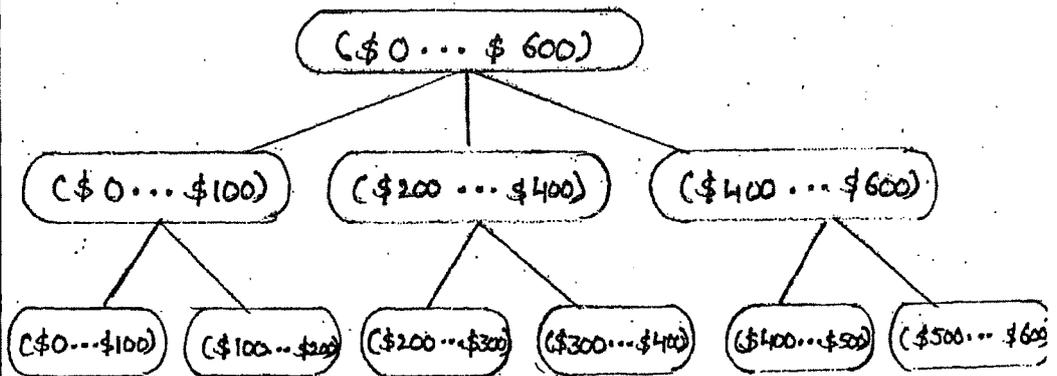


fig (2-2-4-4). Set Grouping hierarchy for price dimension

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2.2.5. OLAP operations in multi Dimensional Data

Model:-

The multi Dimensional Data model allows the data is organized in multiple Dimensions and also Each dimension contains the set of levels as defined in concept hierarchy. The OLAP operations are shown in the following diagram.

combination of data-mart is going to form

DWH

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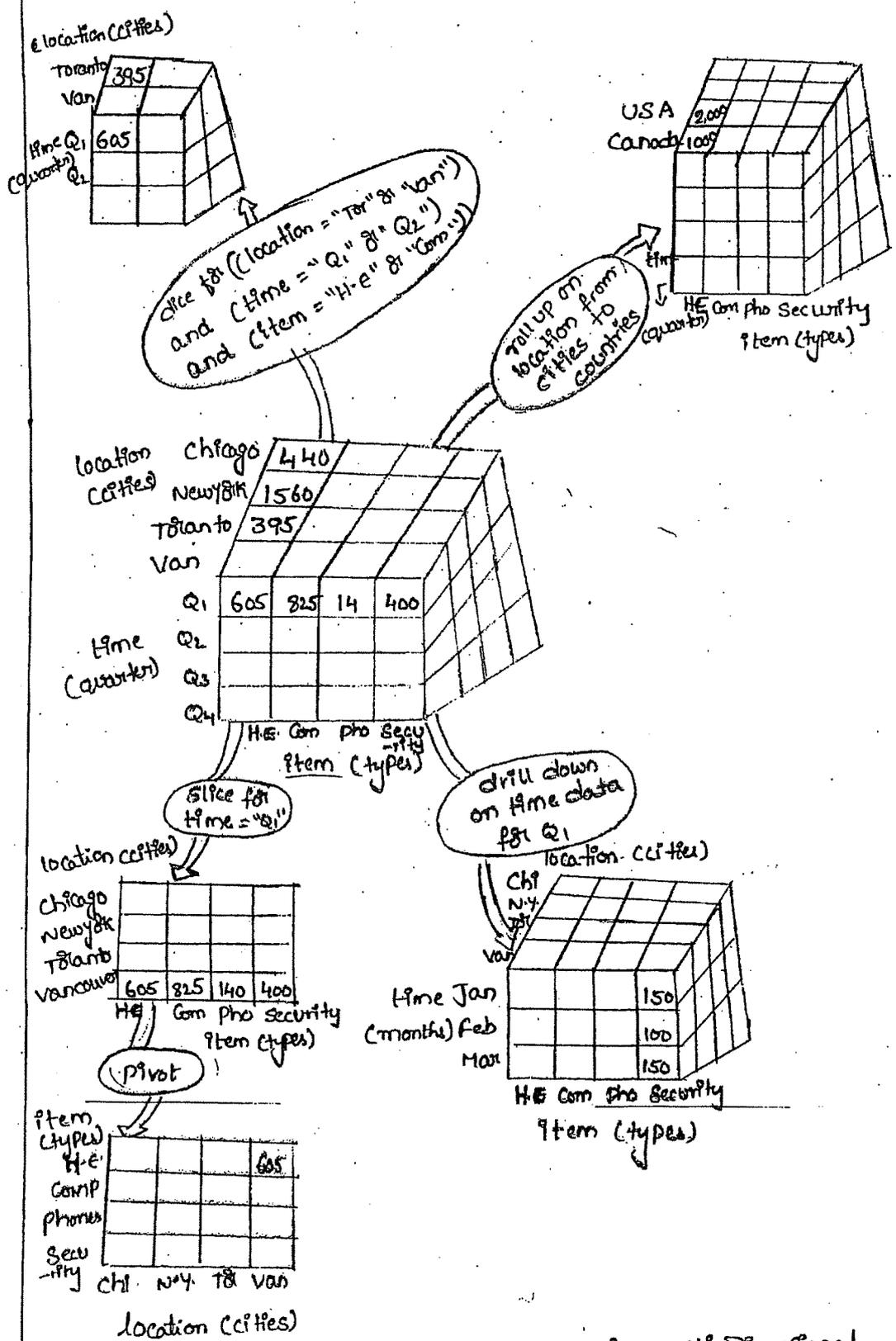


Fig (2.2.5.1): OLAP operations for Multi Dimensional Data Model

(1) Roll-Up:

It is also called as Drill-Up. It is specified by climbing up a concept hierarchy for a given dimension. Here roll-up operation is specified for location (from cities to countries).

(2) Drill-Down: It is the reverse of roll-up operation. Here Drill-Down operation is specified for time data on Q_1 . Drill-Down means it is the stepping down the concept hierarchy for the given dimension.

(3) Slice and Dice: In slice operation we select only one dimension to get the sub cube. Here slice operation is specified for time data = Q_1 .

In Dice operation we select the 2 or more dimensions to create the sub cube. Here Dice operation is specified for location = "Toronto" & "Vancouver", Time = " Q_1 " & " Q_2 ", item = "home entertainment" & "computer".

Pivot: It is the Visualization operation. Here we change the data access to present the another view of data. pivot is also called as "Rotate".

In some of the OLAP applications additionally contains two operations.

1) Drill Across: This uses the relational db requires to access the data from multiple fact tables.

2) Drill-Through: It also uses the relational db queries to drill lowest level of data cube. i.e; down to its back end relational table.

2.2.6. Starlet Query Model for querying multi-dimensional Databases:

The starlet query model consists of radial lines originating from the centre and each radial line represent concept hierarchy for given dimension and also each level in concept hierarchy is called as "foot print".

The starinet query model for All Electronics database.

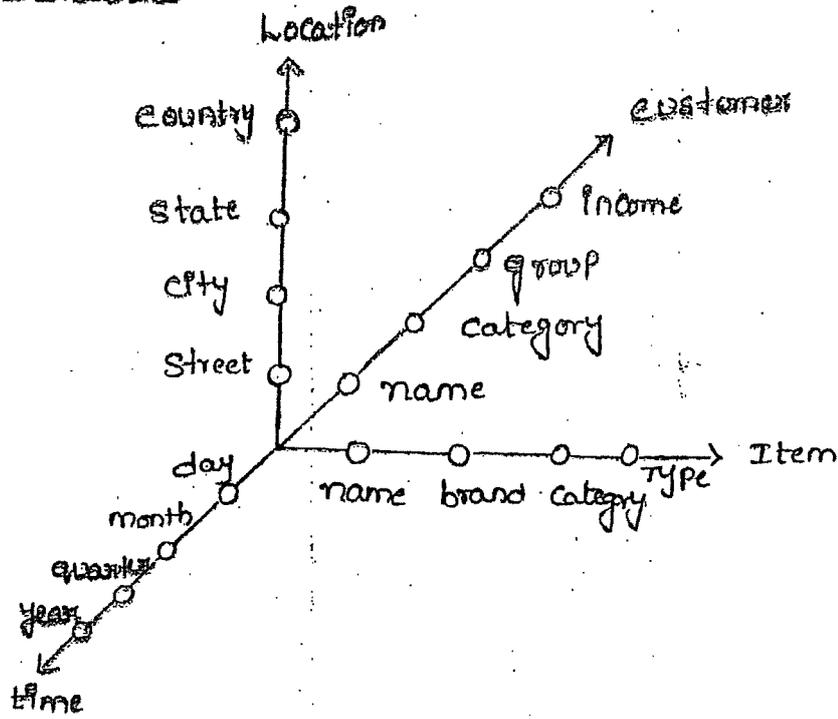


fig (2.2-6.1): Starinet query model for
All Electronics Database

39) 2.3 Data Warehouse Architecture :-

2.3.1. The steps for designing and construction of DWH :-

Reasons for DWH Design :-

1. The DWH provides the competitive advantage.
i.e., it contains the live data.
2. Using DWH we can extend our business.
3. It provides the security.
4. It improves the Customer Relationship.
5. It Reduces the cost by using Relational DB queries to design the DWH generally we follow the 4 views.
 - (a) Top-down view.
 - (b) Data source view.
 - (c) Data warehouse view.
 - (d) Business process view.

(a) Top-down view :-

In this view we select the essential data before going to the DWH design.

(b) Data source view :-

In this view it ~~exposes the~~ data stored & managed by the operational data bases.

(c) DWH view :-

In this view data is used stored in the form of fact & dimensional tables.

d, Business process View :-

Here data is extracted from the DWH and it presents to the end user according to the opinion of the end user.

The process of DWH Design :-

To design the DWH generally we follow the 3 approaches.

1. Top-down Approach.
2. Bottom-up Approach.
3. Combined Approach.

1. Top-down Approach :-

In this approach we start from the strategic planning and design. This approach is best suitable for the technology is known & business problems are solved.

2. Bottom-up Approach :-

In this approach we always perform the experimentation and here the system is completed quickly.

3. Combined Approach :-

In this approach it contains the 2 advantages that is strategic planning from top-down design & Rapid implementation from bottom-up approach.

Generally to design the DWH we follow the any one of the 2 methods. i.e., Water fall Method (a) spiral Method.

Waterfall Model :-

In this model we construct systematic and structured analysis for each step and before going to the next step this is just like the waterfalling from one step to next step.

Spiral Model :-

In this Model updation is easy i.e., it can be adapted to any new technology or any Model. Because of this, it is best suitable for DWH & Data Marts.

The steps for Resigning DWH process :-

The steps are

① choose a business process to Model. i.e., If we require entire data and its scope is enterprise wide then go for the DWH Model.

The Data scope is limited to with in the department then go for the Data Mart Model.

- (2) We have to define the level for each dimension.
- (3) We have to define the dimension tables that must be included in fact Table.
- (4) We have to define the Measures that is to be analyzed like dollars - sold, units - sold.

Three-Tier DWH Architecture :-

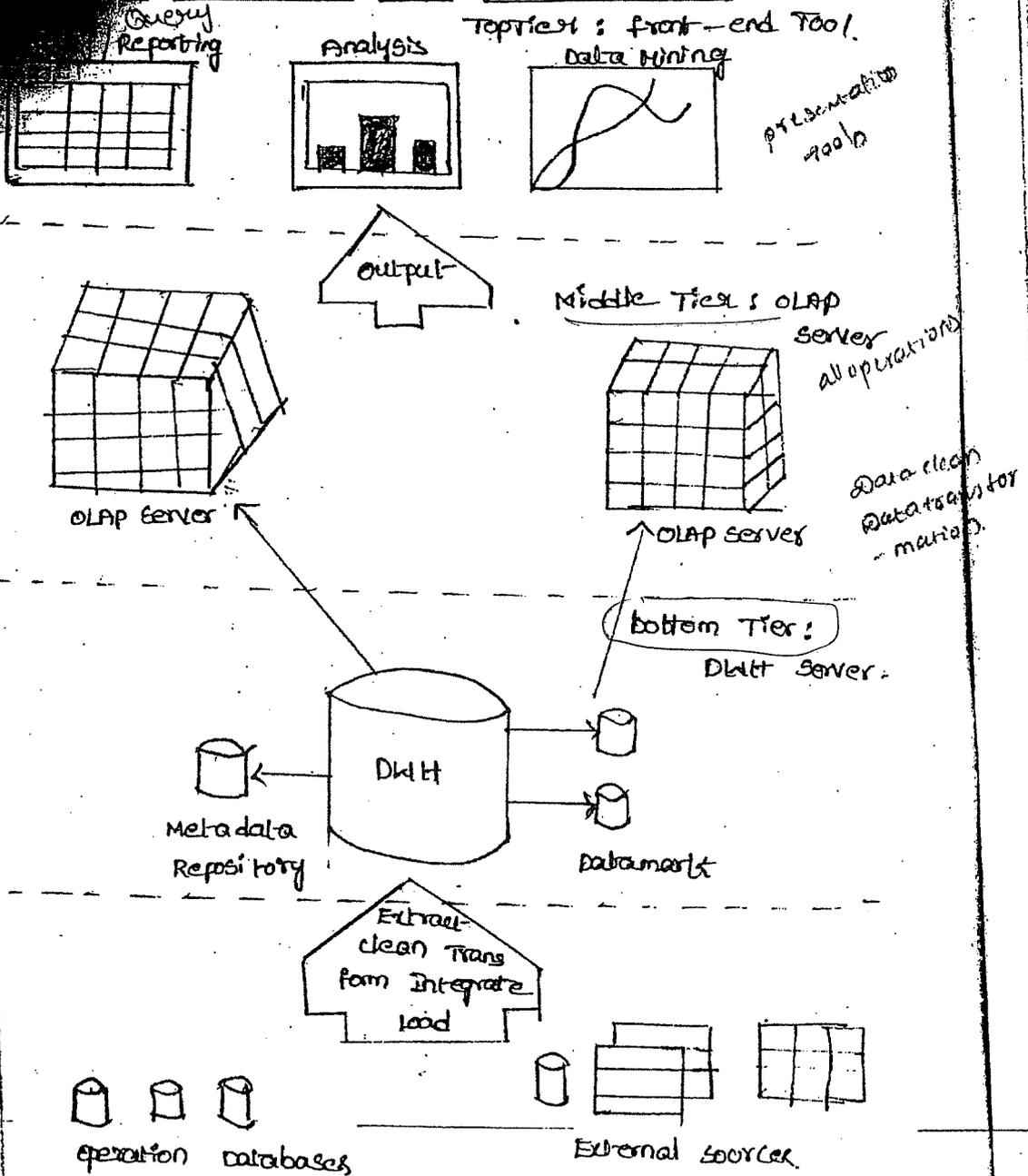


fig (2.3.2) A Three-Tier DWH Architecture.

This architecture contains the 3 Tiers.

1, The Bottom Tier contains the DWH Server - Here data is extracted from the operation databases (or) external sources like Text-file, Excel-sheet, online-data record, etc. Then we apply data cleaning & data Transformation then this entire data is integrated and finally this is loaded into DWH. To extract the data we use the application programs. These application programs are called as gateways.

2, The Middle Tier contains the OLAP servers. These servers are constructed directly from DWH data (or) Data Marts data. To construct these servers we use 2 Models.

- i, ROLAP Model (Relational OLAP) - 3 schemas
star, snowflake, fact constellation
- ii, MOLAP Model (Multi-Dimensional OLAP).

ROLAP Model, it is the Relational db system. It maps the Multi-dimensional data to Relational db systems.

MOLAP Model, These are the special servers. These servers provide the multi-dimensional view.

3, The Top Tier contains the front end tools like query and reporting tools, analysis tools, data mining tools. using these tools the data is presented to management.

Recommended Architecture for DWH

The Recommended architecture contains the 3 steps.

Step 1:

We have to define the high level enterprise data Model. This model must be consist Realible and it must integrate the all the subject areas data.

Step 2:

In step 2, we design independent data marks at the same time we have to develop the enterprise DWH.

Step 3:

In step 3, first of all we develop the distributed data marks. These are used to integrate the different data marks data. Then finally we construct the multi-Tier architecture for DWH.

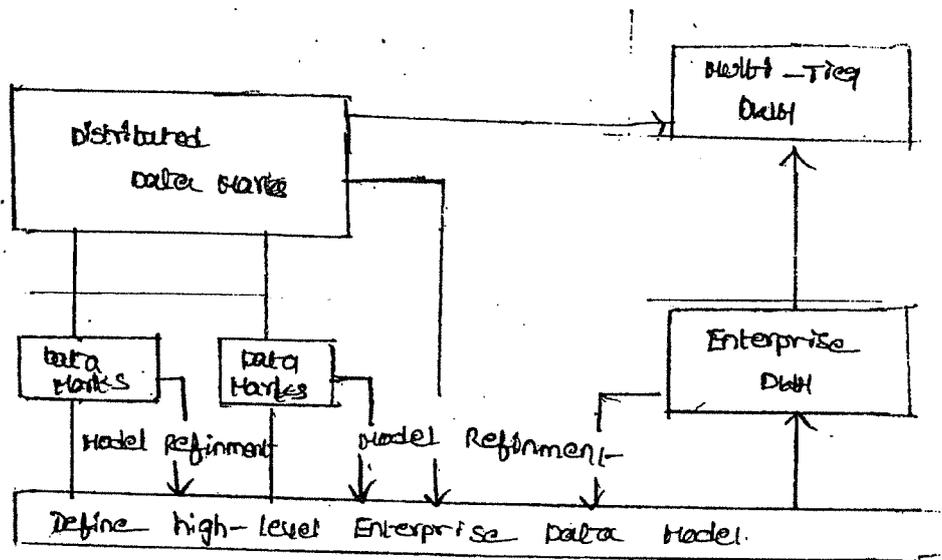


fig (2.3.2.2) A Recommended Architecture for DWH (Two-Tier).

2.3.3 Types of OLAP Servers : ROLAP, MOLAP & HOLAP :-

The types of OLAP servers mainly classified into 3.

1. ROLAP (Relation OLAP).
2. MOLAP (Multidimension OLAP).
3. HOLAP (Hybrid OLAP).

These servers provides the multidimensional view to the end user & the data is accessed from DWH & data marts.

1. ROLAP :- (Relation OLAP)

: This is the extensions of relational db Management system. It is also called as Relational db Management system. It contains the advantage of greater scalability and these ROLAP servers placed in b/w DWH & front end tools.

2. MOLAP (Multidimensional OLAP) :-

These servers store the data in the form of multidimensional arrays and provides the multidimensional view to the end user. These servers contains the advantage of efficient space utilization.

3. HOLAP (Hybrid OLAP) :-

It is the combination of ROLAP & MOLAP. It contains the advantage of scalability from ROLAP & efficient space utilization from MOLAP.

2.4 Implementation of DWH: Using Data cubes and OLAP

The DWH contains the large amount of data. Therefore it is a critical for Data Mining system to provide efficient cube computation techniques & query process techniques.

2.4.1 Efficient computation of data cubes :-

In multidimensional view cube computation extends SQL and includes "compute cube operator". For example sales data cube for all electronics contains 3 dimensions and one measure. i.e., item product year and measure is sales in dollars. This data is analyzed by using the following

- "compute sum of sales, group by item, year"
- ||| "compute sum of sales. group by year"
- "compute sum of sales group by item".

The above cube contains the total no. of cuboids (8) group by's are $= 2^3 = 8$ i.e.,
 { (item, product, year), (item, year), (item, product),
 (product, year), (item), (product), (year), () }

This above cube is represented in the form of lattice of cuboids. This is shown in below.

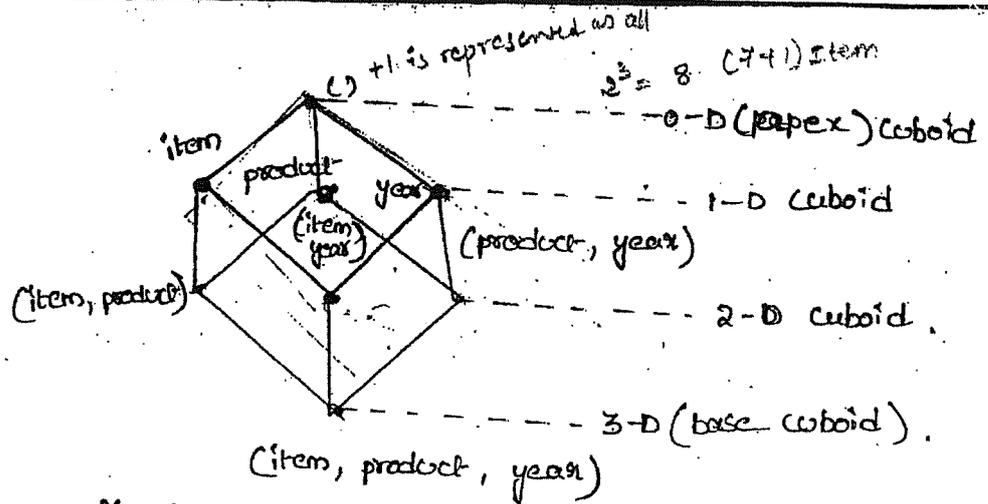


fig (2.4.1): Lattice of cuboids makes the 3-D (database) datacube, the dimensions are item, product, & year.

Compute sales group by item (or) product or year here, apex cuboid is represented by

(1) or all IT Specify group by is Empty total is total sum of all the sales. This is computed by using the query.

"Computed sum of total sales". Similarly one dimension operation is specified by using query

"Compute sum of sales group by item". 2-D operation is specified by using query

"Compute sum of sales group by item and year."

finally 3-D operation is specified by using query.

"Compute sum of sales group by item, product and year".

The above data cube is defined by using the DMQL.

DMQL \rightarrow Data Mining Query Language

Define cube sales {item, product, yearly} :
Sum (sales - in - dollars). A cube with n dimension contains the total cuboids = 2^n , if there is no hierarchy for each dimension. The above contains the 3 dimensions. Therefore total cuboids = $2^3 = 8$. But if the each dimension contains the level of hierarchy like time dimension i.e., day < months < quarters < year; then total cuboids = $\prod_{i=1}^n (l_i + 1)$, where n is the no. of dimensions l_i means levels associated with the dimension i .

Here '1' specifies the high level summarization i.e., represented by R all, or, ()

For example a cube contains 10 dimensions and each dimension contains the 4 levels then approximately. The total no. of cuboids is 5^{10} . Therefore it contains the large no. of cuboids and also each cuboid requires the large amount of space. This is avoided by using

concept hierarchy. - 1 level sales 2 level 3 countries
eg cards \rightarrow countries \rightarrow states

3. The aggregates are computed from the previously computed aggregate rather than taking from fact table.

MOLAP cube computation techniques :-

In MOLAP the data is stored in the form of multidimensional array. here each row is divided into chunk. The chunk is a subcube and small enough to fit in the memory available for cube. The n-dimensional array is divided into n-chunks. This is called as chunking. Each chunk is stored as a object on the disk. Each chunk is identified by "chunkID + offset".

For example 3-dimension array contains 3 dimensions. i.e., A, B & C, then the chunks are shown in below.

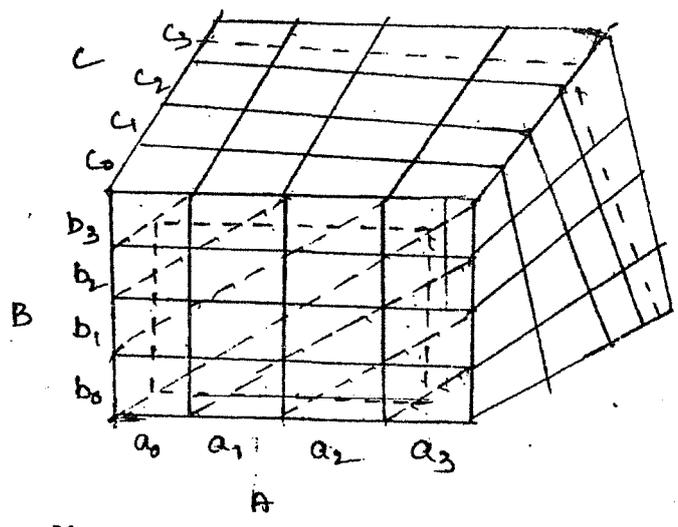


Fig (2.4.2) : 3-D array contains 3 dimensional A, B & C organized into 64 chunks.

partial materialization :-

The data cube materialization mainly classified into 3.

- 1, No Materialization (eg - item)
- 2, Full Materialization
- 3, partial Materialization (2 to 8 dimension & value aggregation)

In those 3rd one is best option i.e., partial materialization. partial materialization means selected cuboids are materialized. i.e., selected cuboids are analyzed. i.e., the selected cuboids data is represented by using data cubes.

The partial materialization contains the 3 steps.

1. Identify the ^{subjects} (subjects) of cuboids that is to be materialized.
2. Analyze cuboids during the query processing
3. update the cuboids during the data load process.

Multiway Array Aggregation in the Computation of Data cubes :-

To Analyze the data ROLAP contains tuple & tables these two are the basic data structures in ROLAP.

ROLAP cube Computation Techniques :

The ROLAP uses the

- 1, Sorting, hashing & grouping operations on dimension attributes to recorder the data.
- 2, The ROLAP uses the grouping for sub queries to partly the processing of sub queries.

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The datacube contains the following.

1. The highlevel summarization is represented by () or all.
2. The 2 dimension i.e., AB, AC & BC is calculated by grouping AB, AC, BC.
3. The one dimension i.e., A, B & C is calculated by grouping dimensions individually.
4. The 3-D (base) cuboid is represented by ABC.

Two arrangements for Multidway array aggregation for 3-D cube concept hierarchy of 3 values & etc.

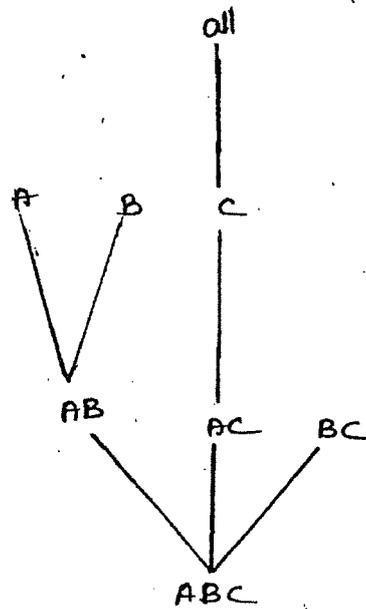


Fig (a) Best used cube computing

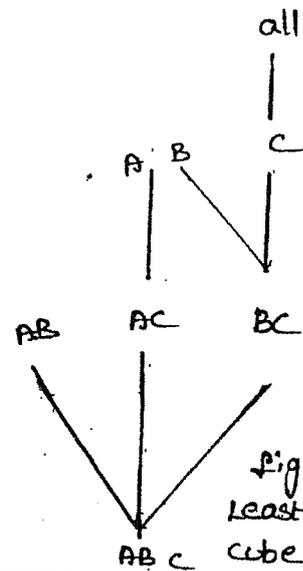


Fig (b) Leastly used cube computing

for ex, the memory units for dimension A, B & C are 40,400 & 4000 memory units. then largest space required for 2-dimension. i.e., BC (for units = $400 \times 4000 = 1600000$ MU) then 2nd largest space is required for 2-D.

i.e., AC (for units = $40 \times 4000 = 160000$ MU). The smallest space is required for 2-D.

i.e., AB (for units = $40 \times 4000 = 160000$ MU).

∴ In fig (a), Minimum space required.

$$= AB \text{ (for whole AB)} + AC \text{ (for one row for AC)} \\ + BC \text{ (for one chunk for BC)}$$

$$= 40 \times 4000 + 10 \times 4000 + 100 \times 1000 \\ = 16000 + 40000 + 100000 = 1,56,000 \text{ MU.}$$

In fig (b), Minimum space required.

$$= BC \text{ (for whole BC)} + AC \text{ (for one row for AC)} \\ + AB \text{ (one chunk for AB)}$$

$$= 400 \times 4000 + \overset{40}{10} \times \overset{1}{1000} + 10 \times 100 \\ = 1600000 + 40000 + 1000 \\ = 16,41,000 \text{ MU.}$$

2.4.2 : Indexing OLAP data :

To provide efficient data access most of the DWH systems contains index structure and materialization. The materialization is provided through data cubes. The index structure is provided by using bit map indexing and join indexing.

Bit map indexing :-

This is the most popular data accessing technique in OLAP. Using this we find the data quickly in data cube. This is the alternative technique for Record-Id (RID) lists. In Bit-Map indexing. Each attribute value

consists of distinct bit vector values. If the attribute value 'V' for the given row in base table then its value is set to '1' corresponding to the row in the bit-map index, all the remaining bits are set to '0' in the corresponding row at bit-map index.

for example consider the all electronics db.

Base Table

RID	item	city
R ₁	H	V
R ₂	C	V
R ₃	P	V
R ₄	S	T
R ₅	H	T

Item Bit-map Index Table

RID	H	C	P	S
R ₁	1	0	0	0
R ₂	0	1	0	0
R ₃	0	0	1	0
R ₄	0	0	0	1
R ₅	1	0	0	0

City Bit-map Index Table

RID	V	T
R ₁	1	0
R ₂	1	0
R ₃	1	0
R ₄	0	1
R ₅	0	1

fig (2.4.2.1): OLAP index by using bit-map indexing

Advantages of Bitmap Indexing :-

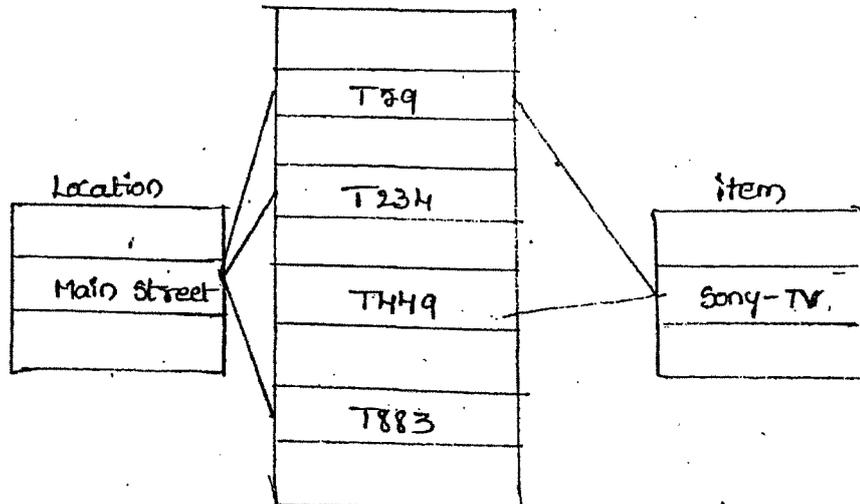
- (1) It is the easy as well as the simple compared with the indexing & hashing.
- (2) It requires the less amount of accessing time because the value is represented by using only 1 bit.
- (3) It requires less amount of space.

Join indexing :-

It is mainly used in relational db by using this we easily find the joinable tuples.

for ex, two relations are A(RID, A) and S(B, SID) join on attributes A and B. The join index record pair (RID, SID) where RID and SID are the record identifiers for relation A and S.

For example consider the sales fact table and 2D tables i.e., location and item. The relation b/w the sales fact table and dimension tables are shown in below.



fig(H.2.2): Relation b/w sales fact table and 2D tables i.e., location and item.

Here, main street value in location dimension joins the tuples with T59, T234 & T883 of the sales fact tables similarly the value Sony TV of item dimension joins the tuples T59 and T449 of the sales fact table.

The join index tables are shown in

below :

join index table
location/sales

location	sales-key
-----	-----
Main-street	T59
Main-street	T234
Main-street	T883
-----	-----

join index table
item/sales

item	sales-item
---	---
Sony-TV	T59
Sony-TV	T449
---	---

Join Index Table for Two Dimensions
Location / Item / Sales.

Location	Item	Sales-Key
Main-street	Sony TV	T59

Fig (2.4.2.3) Join Index Tables.

This join index Tables mainly used to maintain the relation b/w primary key and foreign key. This is mainly used in star schema to maintain the relation b/w foreign key of the fact table with primary key of dimension tables.

2.4.3 Metadata Repository :-

Metadata is nothing but data about the data. The Metadata repository contains the following steps.

1. The structure of DWH which includes schemas, views, dimensions, hierarchies and data definitions.
2. The operational repository which includes data usage and monitoring information.
3. It also includes the algorithms for summarization.
4. To map data sources to data warehouse which includes data cleaning, data transformation, data integration and security techniques.
5. The business repository which includes business policy, conditions, terms and business definitions.

2.5 Datacube Technology :-

Here initially, we write about the data cube in multidimensional data model.

It is mainly classified into.

2.5.1 "Discovery Driven Exploration" of data cubes :-

Here analyst or user search for the interest-patterns by using OLAP operations i.e., drill-down, drill-up, slice and dice & pivot. OLAP operation imp

If the discovery process is not automated then end user search for the interesting patterns manually this is difficult. This is avoided by using the method discovery driven exploration. In this discovery driven exploration we identify the interesting pattern and it is marked. To provide the visualization to the end user.

Discovery Driven Exploration :-

In this method previously computed measures indicates the data exceptions after that these measures treated as exception indicators and using this end user identifies the interesting pattern easily.

∴ these measures provide the "degree of surprise" i.e., ratio of existing value in the cell with expected value.

These measures are calculated by using statistical analysis or grouping functions. These measures mainly classified into 3.

Chapter-3

Data Preprocessing

Database consists of massive volume of data which is collected from heterogeneous sources due to this heterogeneity, real world data tends to be inconsistent and noisy. If data is inconsistent, then there is a possibility that mining process can lead to confusion which results in inaccurate.

Need for preprocessing the Data;

Incomplete, noisy and inconsistent data is common in large real world databases and datawarehouse

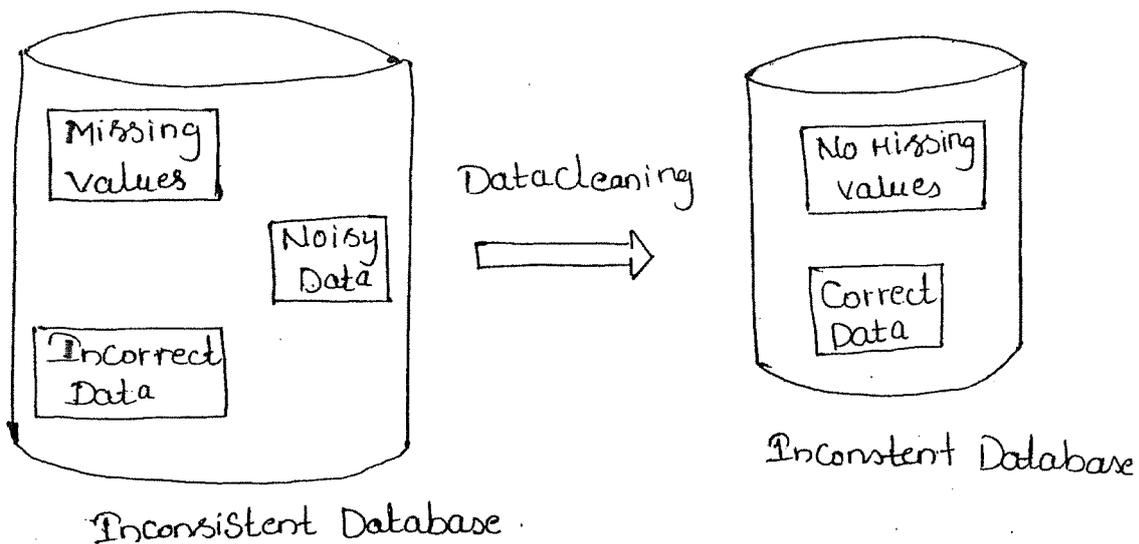
Incompleted data can occur for a no. of reasons.

1. Attributes of interest may not always be available.
2. Relevant data may not be recorded due to a misunderstanding.
3. Data that is inconsistent with other recorded data might be deleted.
4. The data collection instruments used may be faulty.
5. There may have been human or computer errors occurring at data entry.
6. Errors in data transmission can also occur.

To overcome the above problems the following data preprocessing techniques are required.

1. Data cleaning
2. Data Integration
3. Data Transformation
4. Data Reduction.
5. Data Discretization.

1. Data Cleaning: when data is collected from data sources then there are chances that the data can be inconsistent, incomplete and noisy. Data cleaning is a process of removing unnecessary and inconsistent data from the databases. The main purpose of data cleaning is to improve the quality of data by filling missing values, reconfiguring the data to make sure that data is in consistent format.



Fig(1): Data cleaning

1.1: Missing Values:

The missing values consists the following techniques and corrected before applying DMQL.

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a) Ignore the tuple:

This technique is simple. In this just we avoid the tuple which doesn't contain values. But it's not recommended.

b) Fill the missing values manually:

Here we manually fill the missing values. But this technique is not recommended for large database which contains important values.

c) Using Global Constant values:

Here the missing values has been filled with the global constant such as ^{as} unknown as (∞) infinity. But this technique not succeeded because it deviates the Data mining Process.

d) Using attribute mean value:

Here each missing value has been filled by mean value i.e. In All Electronics customer average income is 2800 \$. Then the customer record i.e. the income attribute's missing value is filled with 2800 \$.

e) Use the most portable values:

Here we find out the most portable values by using different techniques like Bayesian classification and decision tree induction etc.

1.2 Noise data :

Here the noise data has been smoothing out by comparing neighbouring values. In this technique

data values has been distributed into different wings.

1. Binning Method :

a) : Smoothing by Bin-by-mean

Here we find out the mean value for each Bin & missing values are replaced with that mean value.

b) : Smoothing by Bin Bounding :

Here we find out min and max values for each Bin then each Bin's value is replaced with the closest value to the min or max.

Ex: price in dollars of an item is as follows.

4, 8, 15, 21, 21, 24, 25, 28, 34.

Bin 1 : 4, 8, 15

Bin 2 : 21, 21, 24

Bin 3 : 25, 28, 34

Smoothing by Bin

Bin 1 : 9 9 9

Bin 2 : 22 22 22

Bin 3 : 29 29 29

Smoothing by Bin Boundary value;

Bin 1 : 4 4 15

Bin 2 : 21 21 24

Bin 3 : 25 25 34

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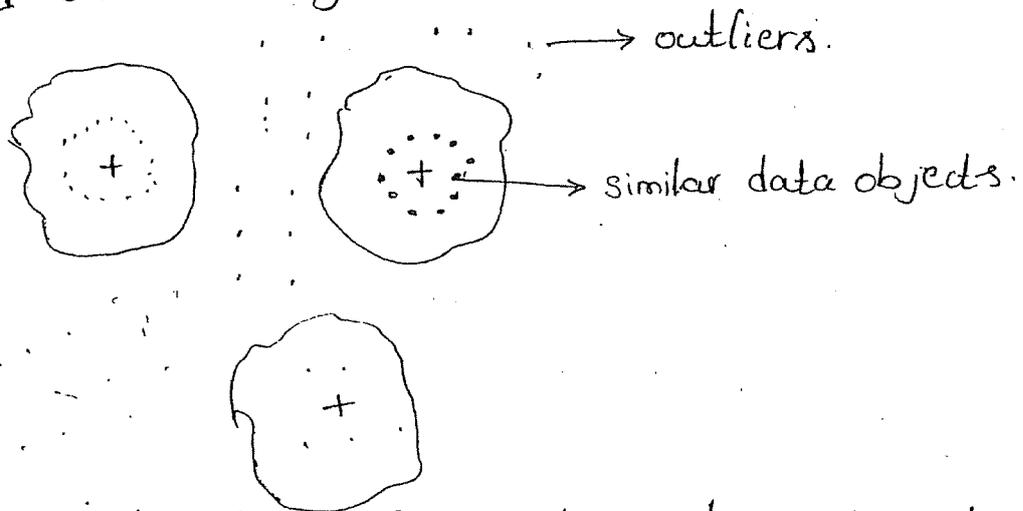
ings.

In Bin boundary we identify the data if it is near to the min value then we replace that value with min I or min values otherwise if it is near to the max value then we replace that value with the max value.

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Clustering: In this we find out outliers i.e. clustering identifies similar data objects those are placed in one cluster & also identifies dis-similar data objects and those are called outliers, in other words grouping the similar data objects into 1 place is called as clustering.



Using combination of computers & human inspection:

In this outliers can be identified through a combination of computers & human inspection i.e. we need to identify any algorithmic approach to find out the outliers and clusters in our data set along with this we have to take the help of manual procedure to identify clusters & outliers.

Inconsistent data:

Inconsistent data means the data with

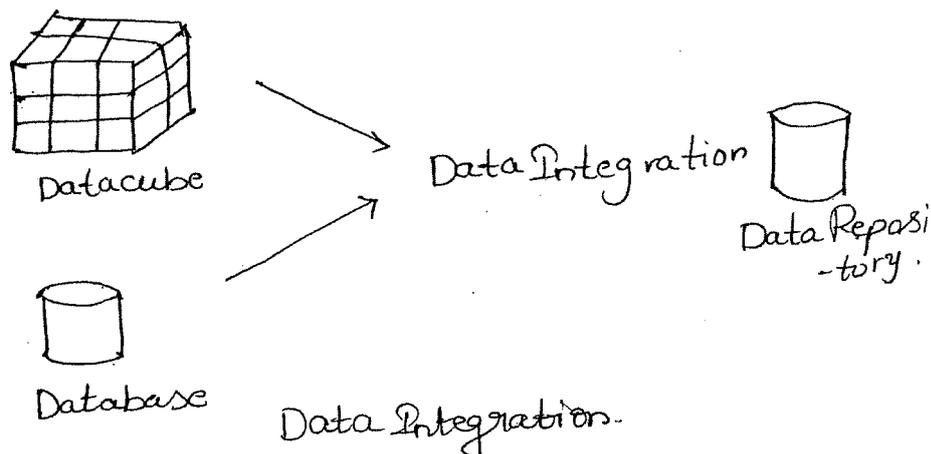
duplicated values. For Eg; ItemId is used to categorize items in AllElectronics. If the item-id entered wrong or missing then we can say that item-id is inconsistent data. Then by applying different smoothing techniques Bin by mean, Bin by Boundary etc. to convert the data into consistent data.

Data Integration & Data Transformation:

Here the data from different data sources is collected and placed into the DWH, but it contains different difficulties to make unique identification or unique representation for the data of different data sources.

Data mining requires data integration. Data integration is the merging of data from multiple data stores into a coherent data store as in DWH. These sources may include multiple databases, datacubes or flat files.

Data Integration: Data integration is a process of combining data from heterogeneous data sources such as different data bases, flatfiles etc, to form a single consistent data repository.



For eg: In one data source customer-id is entered as cust-id where as in another data source it entered as Customer Number.

The integration contains different problems while collecting the data from multiple tables. This can be avoided with an analysis called as correlation analysis. For Eg: The correlation analysis or relation b/w two attributes can be measured as,

$$= \frac{\sum (A - \bar{A})(B - \bar{B})}{(n-1) \bar{\sigma}_A \bar{\sigma}_B}$$

In the above Equation 'n' represents no. of tuples \bar{A} & \bar{B} represents mean values of A, B $\sum A$ & $\sum B$ represents standard deviations of A, B.

we can find the mean value of A as $\bar{A} = \frac{\sum A}{n}$

Similarly we can find out the standard deviation of A as

$$\bar{\sigma}_A = \sqrt{\frac{\sum (A - \bar{A})^2}{n-1}}$$

If the result of correlation analysis is greater than 0 (zero) then two attributes are positively correlated i.e., if we increase the value of attribute A then the value of attribute B will also increase.

If the correlation analysis result is '0' then two attributes are independent to each other.

If the correlation analysis result is less than 0 then the 2 attributes are negatively correlated i.e. If the value of A increases then the value of B will decrease.

Data Transformation:

Data Transformation is nothing but converting different data sources into a format that must be acceptable for the data mining system.

5, 48, 99, 35, 81 Data Transformation 0.005, 0.048, 0.049, 0.035
0.081.

1. Smoothing:

Here the noise data is subjected for smoothing by using different smoothing techniques, Binning, clustering etc.

2. Aggregation:

Here we use some aggregated functions to perform data transformation. For eg; using daily sales we can compute monthly sales. In the same way, using monthly sales we can compute yearly or Annual sales.

3. Generalization:

Here lower level value is replaced with higher level value by using concept hierarchy. For eg: The dimension location contains concept hierarchy street, city, state & country instead of placing the street value would be more generalized to place Country value; the data mining system.

4. Normalization:

In Normalization we force each value should in specific range. -1.0 to +1.0

5. Attribute Construction: In this new attribute

will be constructed by integrating attributes from different sources.

In the above 5 methods the Best one is

"Normalization".

Normalization;

It contains following techniques.

1. Min-Max Normalization:

Here \min_A \max_A are the minimum & maximum values of the attribute A. Then the value v of A is transformed into v' . By using new range of new \min_A & new \max_A .

we compute new values i.e. v' By using below formulae.

$$v' = \frac{v - \min_A}{\max_A - \min_A} (\text{new } \max_A - \text{new } \min_A) + \text{new } \min_A$$

min max Normalization maintains same relationship of original data values. For Eg: The income attribute contains the min & max values of \$ 12000 & \$ 98000 then the income attribute changed into new range (0,1).

By using min, using min max Normalization the original value is \$ 73600 will be transformed, into

$$v' = \frac{73600 - 12000}{98000 - 12000} (1 - 0) + 0.$$

$$v' = 0.716$$

2. Z-Score Normalization:

Z stands for zero mean. so this normalization is also called as zero mean normalization. In this attribute is find out by using mean & standard deviation value.

The value v of A will be transformed by using the formulae.

$$v' = \frac{v - \bar{A}}{\sigma_A}$$

Here \bar{A} is called as mean value σ_A is called as standard deviation. for Eg: The attribute 'A' mean & standard deviation values are \$ 54000 \$ 16,000 then we can Z-Score the Normalization F_i value v \$ 73600 will transformed into

$$\begin{aligned} v' &= \frac{73600 - 54000}{16000} \\ &= 1.225 \end{aligned}$$

Normalization by decimal scaling:

In this technique we move the decimal value based on absolute value of an attribute. By using this technique the original B of A is transformed into B' by using.

$$v' = \frac{v}{10^j}$$

where j is the smallest integer based on the attribute value.

for Eg The attribute A contains values between -986 to +917. Then we find out the absolute value as 986 by using the decimal scaling we can find out the new value v' by dividing it with 1000
i.e $j=3$

$$J=3$$

$$v' = \frac{-986}{10^3}$$

$$= -0.986$$

Data Reduction:

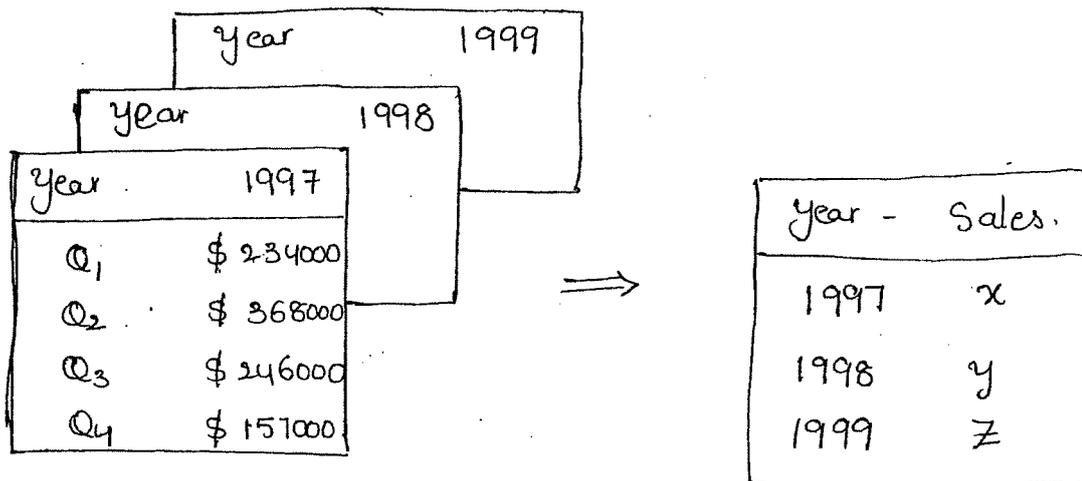
Here we apply aggregation operations, Redundant attributes are removed from the data. Then we can apply data compression technique on data.

In data reduction, we reduce the size of the data in such a way that essential features does not effect. It contains several techniques.

i) Data aggregation:

In Data aggregation we will apply some aggregation operation on the data to construct the datacube & we can also identify some clustered data to reduce the size of the data.

For eg, In ALL Electronics database, the data is stored quarterly based for 3 years i.e, for the year 1997, 1998 & 1999 which is shown in the below diagram.



We can apply different aggregation operations for the annual sales to represent the data in yearly format rather than storing it into quarterly wise.

In the reduction, we can also represent the data in data cube format to represent multiple dimension which is represented in the below diagram.

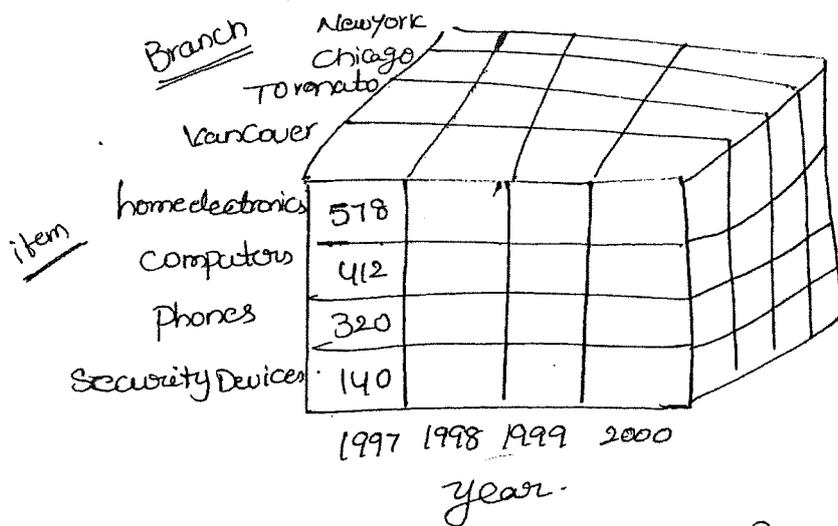


fig: Sales datacube for all electronics.

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ii) Dimensions Reduction: Here the redundant attribute or dimension is identified & it is removed. It contains several methods to reduce dimensions. In those methods, the main aim is we need to identify the min no. of attributes to represent the dataset. Among the available techniques, the greedy method is important. In this method, it enables us to find out the best attribute or worst attribute based on decision tree analysis. It contains the following techniques.

1. Forward Selection: For Eg: Consider the following decision tree.

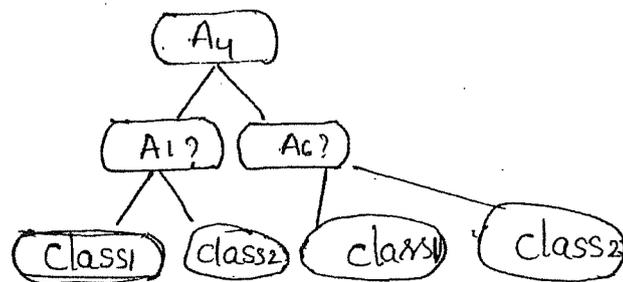


Fig: Decision Tree data.

In forward selection we will start the identification with the empty set. Consider the data set as $\{A_1, A_2, A_3, A_4, A_5, A_6\}$ initial set is $\{\}$ (Empty Set). In the order the first attribute to consider these $\{A_1\}$. The next dataset we consider is $\{A_1, A_4\}$ & the next data set we consider is $\{A_1, A_4, A_6\}$ we leave other attributes which are not important.

2. The Backward Eliminations:

Here we start our searching from the complete

Dataset. In each step we eliminate worst attribute & that is removed from the dataset. This process will be continued until all the dataset is complete. Here the dataset is $\{A_1, A_2, A_3, A_4, A_5, A_6\}$.

So we start over search from this only. In first step, we eliminate the attribute $\{A_2\}$ so the dataset becomes.

$\{A_1, A_3, A_4, A_5, A_6\}$.

\Downarrow
 $\{A_1, A_4, A_5, A_6\}$

\Downarrow
 $\{A_1, A_4, A_6\}$.

3. Combination of forward selection & Backward elimination: It is the integration of above 2 techniques. This is basically applicable for large Dataset. In this we need to identify the best attribute as well as worst attribute based on our requirement. we will add the best attribute to the dataset at the same time we remove the ~~A~~ worst attribute from the Dataset.

To perform all these redundant techniques we need to use decision tree induction. In decision tree induction, we perform test on attributes & we identify the results based on the classifications. Each classification is representative^{ed} by "class labels". In the last example, the total dataset i.e. $[A_1, A_2, A_3, A_4, A_5, A_6]$ is reduced as $[A_1, A_4, A_6]$ & final result

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will fall into either class 1 or class 2.

Definition of Data Reduction :

Data reduction is a process of compressing massive volume of data into limited data set without sacrificing data integrity.

Year 1998	
Sem	Total
1	850
2	800

Year 1999	
Sem	Total
1	825
2	830

Year 2000	
Sem	Total
1	800
2	800

Data Reduction
→

Year	Total
1998	1650
1999	1655
2000	1600

iii Data Compression

It means we reduce the data in such a way that it must be in smaller size i.e., we reconstruct the data from the data in the form of compressed. If we compress the data without any loss of information that Data Compression technique is called as "Loss Less Data Compression" otherwise it is called as "Loss Data Compression".

This Data compression contains two techniques

1. Discrete wavelet transform (DWT)
2. Principle Components Analysis (PCA).

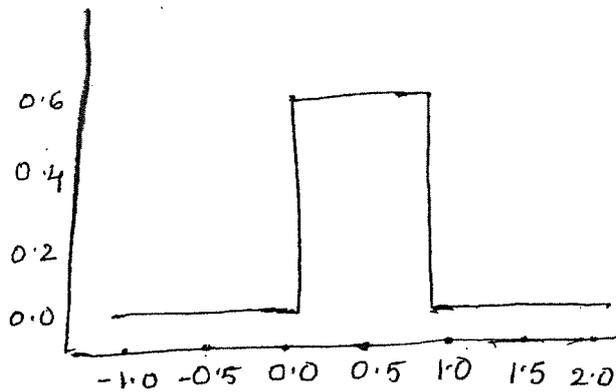
1. Discrete wavelet transform :

In this technique, the data vector D is transformed into D' of wavelet co-efficient. In this technique, we transform any value into new value by considering the coefficients with specific Range. This technique is derived from DFT (Discrete Fourier Transform). This DWT contains several wavelet Transform techniques. But familiar ones are.

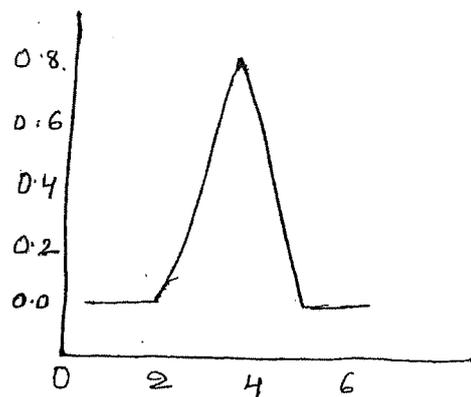
(a) Haar-2

(b) Daubechies 4

which are represented in the below diagram.



(a) Haar-2



(b) Daubechies-4

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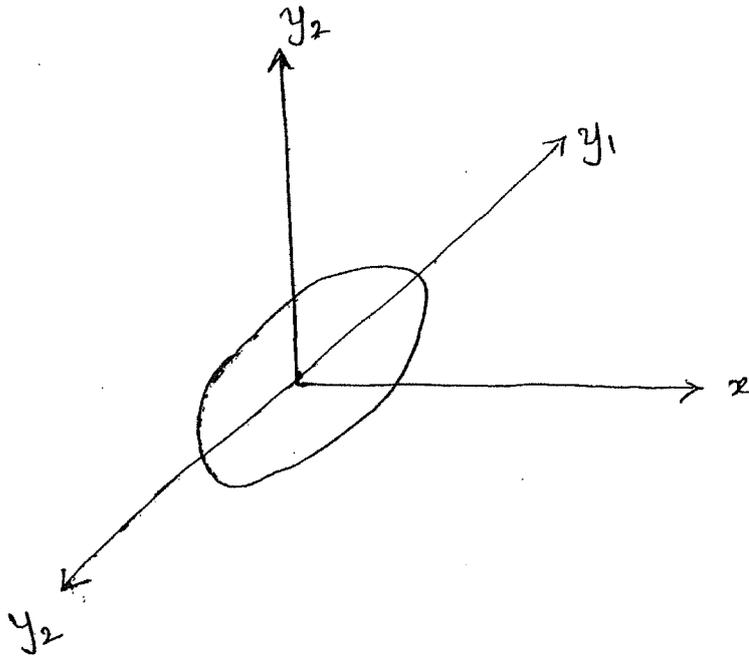
In Haar-2 technique we will consider the data based on the specific range co-efficient r . In the above diagram, the range for considerable values is from 0 to r . Then we consider the class labels in that specific range by leaving other one.

Daubechies 4 also follows same technique but represent in the form of pyramid. The only difference is in Daubechies 4, the dataset length L is represented in the form of even values or in the form of L integer power of 2. Here we apply two functions to compress the data. In that initially we apply smoothing techniques such as sum & average and we find out the differences. If the dataset is too large, then we can divide the dataset consider its length as $L/2$.

The procedure will repeat until the length of the dataset becomes 2. Then we will get compressed data which can be placed into the Data Mining System.

2. Principle Components Analysis:

Here we compress the data of N -tuples or N Data Vectors reduced into K -dimensional orthogonal vectors. By using this technique we can represent the data into compressed vectors. This technique is called as KL (Karhunen-Loève) technique.



In principle component of analysis we apply the normalization of the dataset to represent the entire values in specific range we identify k -dimensional orthogonal vectors & then we compute orthogonal vectors C with new range such that $C < = k$.

Here we construct new access to represents the compressed dataset. According to the above diagram x_1 & x_2 are original axis and y_1 & y_2 are new axis for the original data which represents the compressed data from the original data.

In this method, final higher principle components are integrated to represent the original data.

iv Numerosity Reduction :

Here we reduce the data by selecting the smaller group of values. It is classified into.

- i) parametric
- ii) Nonparametric

Parametric : In parametric Reduction technique, the data is reduced by using

- i, Linear regression
- ii, Log-Linear regression models.

→ In Linear Regression model, we find the best line to be fit b/w the pair of attributes. If 1 value is given then we find the another value. Here we use the formula.

$$y = \alpha + \beta x$$

where α, β are the regression coefficients.

x is the predictor value and

y is the Responsive value.

→ In Log Linear model we predict the cell value of base cuboid. Using this we find the high level cuboids values. i.e By using low level cuboid values we can find high level cuboid values.

Non-parametric:

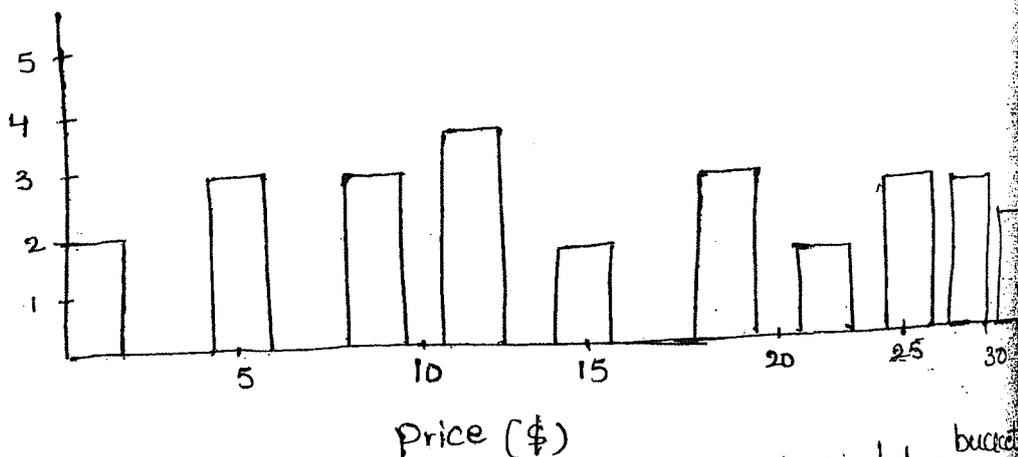
Here data is reduced by using the several methods. The methods are.

1. Histograms
2. Clustering
3. Sampling

1. Histograms:

Here we use the several smoothing techniques. The histogram of an attribute 'A' is partitioned into disjoint parts. These are called as buckets. These buckets are represented in horizontal axis and the bucket count i.e., frequency is represented in vertical axis.

For example, consider the all electronics sales data for price is specified in dollars i.e.; 1, 1, 5, 5, 8, 8, 8, 10, 10, 10, 10, 15, 15, 18, 18, 18, 21, 21, 25, 25, 25, 28, 28, 28, 30, 30. Then the histogram for the above data is show in below.



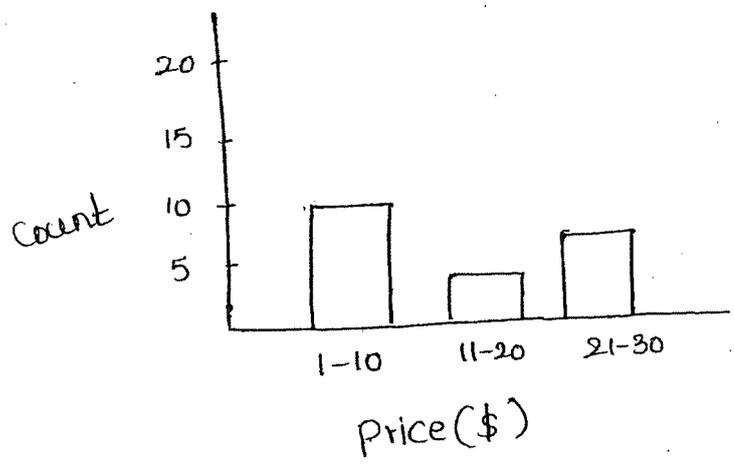
Bucket: (15) number of price according to singleton bucket

L

Here, each bucket contains only one value then those buckets are called as singleton buckets to partition the data we use several partitioning techniques. In those familiar one's are;

i. Equi width;

The equiwidth histogram contains equal width² or constant width. The equiwidth histogram for the above data.



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sales

5, 5,

1, 25,

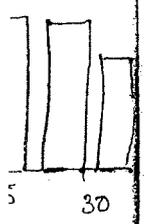
the

ii. Equidepth

The equidepth histogram contains the constant frequency that is, each bucket approximately contains the equal no. of samples.

iii) v-optimal;

The v-optimal histogram contains the low variance. The histogram variance is nothing but sum of values of the bucket.



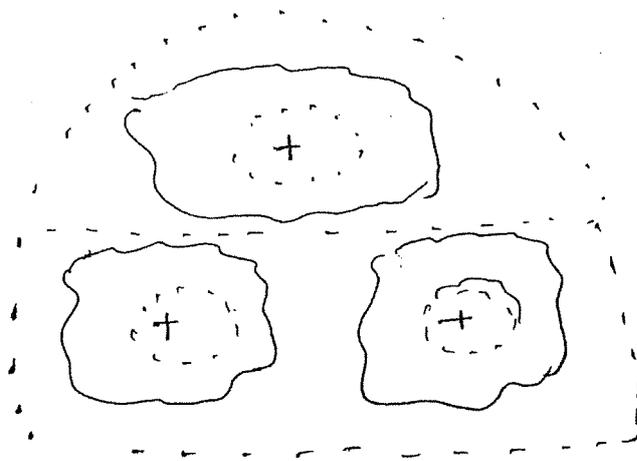
buckets on x

iv) Max-Diff:

The Max-Diff histogram contains the maximum difference of each pair of adjacent values. Here, we use the formula " $\beta - 1$ ", Here ' β ' is the maximum difference. This is specified by the end-user.

2. Clustering:

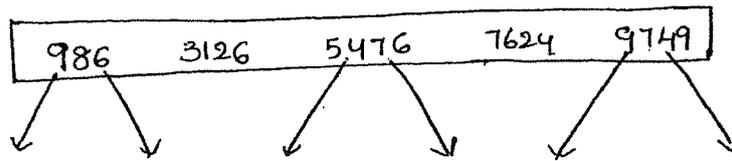
In clustering similar data objects are stored in one cluster and dissimilar data objects are stored in another cluster. For example, 20 customer data w.r. to customer locations in city is shown in below.



Fig(a): 20 customer Data w.r. to customer location in city.

Here center of the cluster is marked with '+'. But in data reduction requires the original values to store the data in efficient way and access the data in efficient way. We use multidimensional index Tree structure. B⁺ tree organization.

The B⁺ Tree organization is shown in below.



Fig(b): B⁺ Tree organization for data set.

For ex consider the large database. This database contains the 10,000 tuples. These are represented by using the keys 1 to 9999. We partitioned this data by using the equidepth histogram. i.e. Each bucket approximately contains the equal no. of values. Then we get the 5 keys. These are ranging from 1 to 985, 986 to 3125, 3126 to 5475, 5476 to 7623, 7624 to 9748, 9749 to 9999.

Therefore, Each bucket approximately contains the $10,000/6$ and also the keys are again divided it into subkeys.

3. Sampling:

In sampling we reduce the data i.e. Large amount of data is replaced with small random sample.

For eg. the dataset "D" and it contains the 'N' no. of tuples. Then the samplings are:

1) Simple Random Sample without Replacement (SRSWR)

for size n:

Here we select the 'n' tuples where $n < N$. The probability of drawing the tuple from D is $1/N$. Here we draw the tuple but, that tuple never be rewritten.

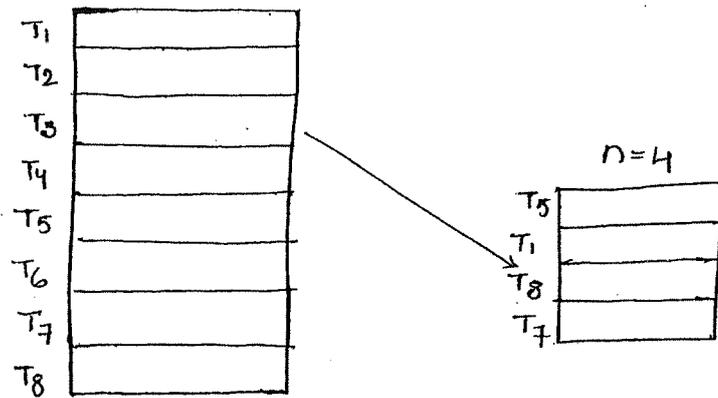


fig (a) : SRSWOR for $n=4$

ii) Simple Random sample with Replacement (SRSWR)
 for size n :

Here, we access the tuple, recorded and rewritten. This technique allows the end user. The tuple a is accessed again.

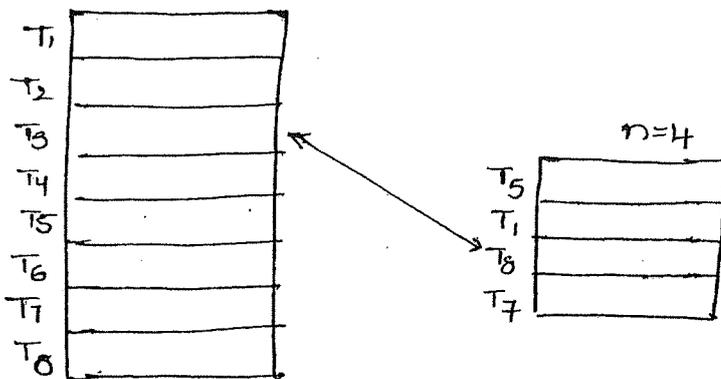
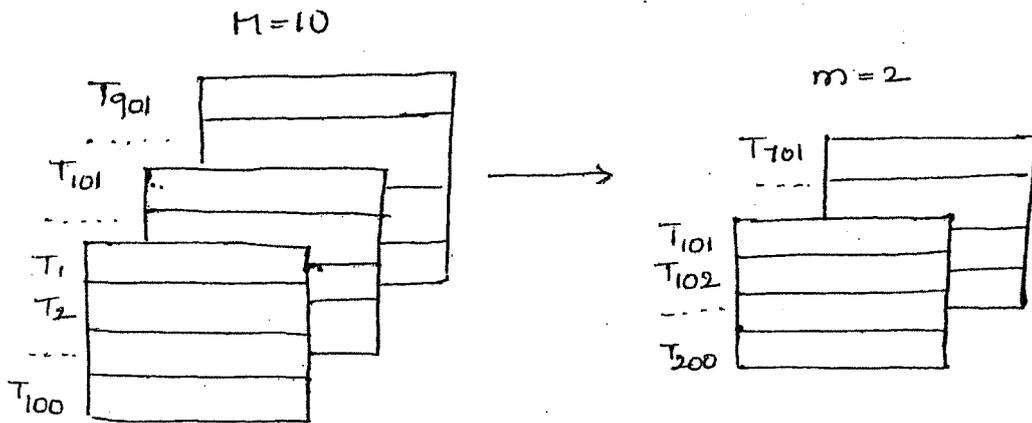


fig (b) : SRSWR for $n=4$.

iii) cluster sample:

Here the tuples in 'D' grouped into 'H' mutually disjoint clusters. Then we select 'm' clusters. This is shown in below.



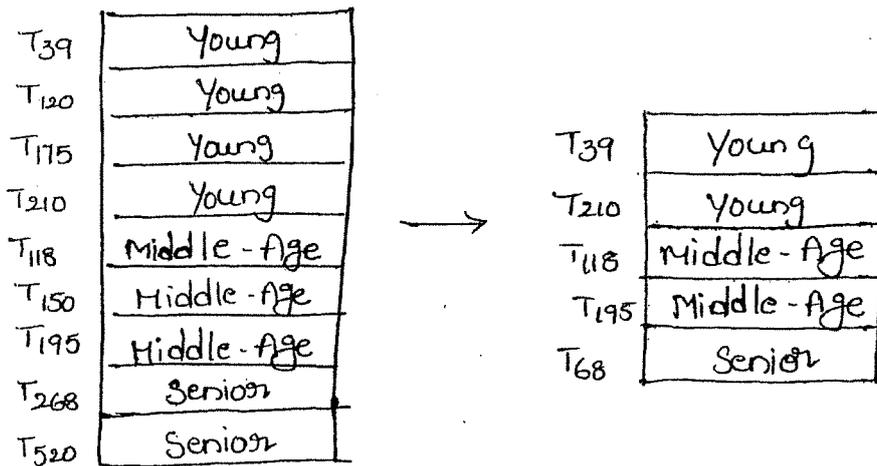
fig(c): cluster sample with $m=2$.

(R).

iv) Stratified sample;

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is

Here, large tuples of 'D' partitioned into mutually disjoint parts. These parts are called as strata. Then we generate the stratified samples of D for each stratum. This is shown in below.



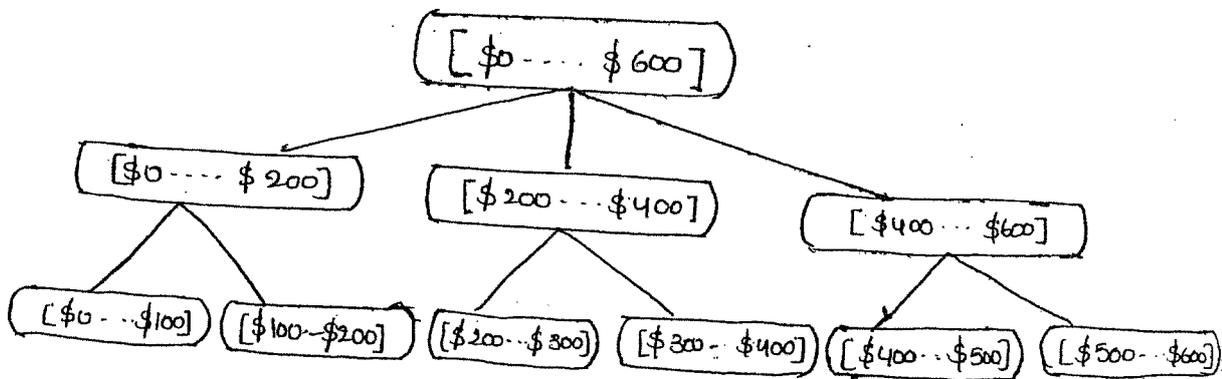
fig(d): stratified sample according to age.

Discretization & concept hierarchy Generation;

The Discretization means we reduce the data value of an continuous attribute by dividing it into intervals. In this best method is concept hierarchy. In concept hierarchy we collect and replaced with

low level hierarchies. For ex. Age attribute contains the values young, middle and seniors.

Then the Age attribute is replaced with one of these 3 values. Concept hierarchy for price is shown in below.



fig(a): Concept hierarchy for price.

Discretization and Concept Hierarchy Generation for Numeric data:

We apply the concept hierarchy easily for Numeric data by using the method distributed attribute analysis. This method contains the 5 techniques.

1. Binning
2. Histogram Analysis
3. Cluster Analysis
4. Entropy - Based Discretization
5. Data Segmentation by natural partitioning

1. Binning:

Here data is partitioned and distributed it into different buckets or bins. Here we use the several Smoothing techniques.

Smoothing by bin Mean:

Here, we find the mean value for each bin and

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that is replaced with each value in bin.

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ii. Smoothing by bin Boundary:

Here we find the maximum value, minimum value. Then each bin value is replaced with closest boundary for ex, Consider the price in dollars are 4, 8, 15, 21, 21, 24, 25, 28, 34. Partitioned with equidepth i.e., 3.

bin 1: 4 8 15

bin 2: 21 21 24

bin 3: 25 28 34

sc - \$600

Smoothing by bin Mean:

bin 1: 9 9 9

bin 2: 22 22 22

bin 3: 29 29 29

for

Smoothing by bin boundary:

bin 1: 4 4 15

bin 2: 21 21 24

bin 3: 25 25 34

analysis

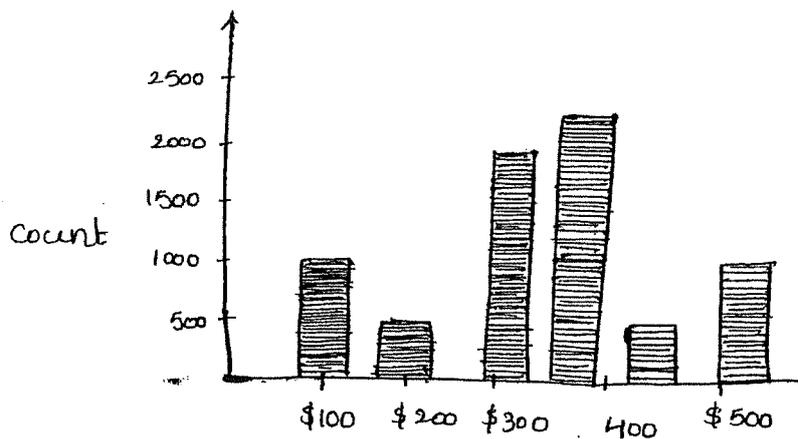
2. Histogram Analysis:

The Histogram for an attribute A is composed by partitioning data into disjoint bucket. These buckets are represented in horizontal axis and the bucket count i.e. frequency is represented in vertical axis. For ex, consider the most frequency range for price is \$ 300 to \$ 350 and this data is partitioned by using Equiwidth partition i.e. Each bucket contains the

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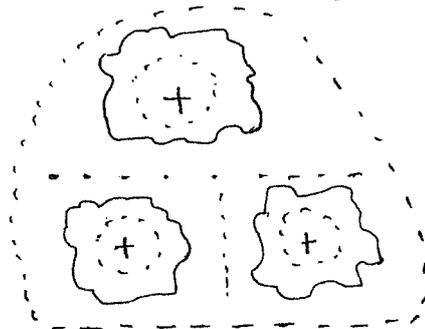
Equiwidth



fig(a): Histogram Analysis for price (\$).

3. Cluster analysis:

In cluster analysis, similar data objects are stored in one cluster and dissimilar data objects are stored in another cluster. The 2D customer data with respect to customer locations in city is shown in below.



fig(a): 2D customer data w.r to customer locations in city. Here center of the cluster is marked with '+' and the cluster again divided it into smaller clusters.

4. Entropy Based Discretization:

The information based measure is called as Entropy. This is applied recursively for the numeric data of Attribute A.

Let us consider 'm' classes. The sample 'S' contains s_i samples in class C_i for $i=1, 2, 3, \dots, m$. Then expected

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information for the given sample.

$$P(s_1, \dots, s_m) = - \sum_{i=1}^m \frac{s_i}{s} \log_2 \frac{s_i}{s} \rightarrow \textcircled{1}$$

Then we can calculate Entropy for an attribute A as

$$E_A = \sum_{i=1}^s \frac{(s_{i1} + \dots + s_{mi})}{s} \times P(s_{i1}, \dots, s_{mi}) \rightarrow \textcircled{2}$$

Finally we can calculate information gain as follows

i).

$$\text{gain}(A) = P(s_1, \dots, s_m) - E(A)$$

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Based on the information gain, we can identify strongly relevant attributes and weakly relevant attributes.

5. Data segmentation by Natural partitioning:

Many of the users prefer the data is distributed uniformly. The uniform distribution allows to the end-user. The end-user read the data easily. For ex, Annual Salaries of a particular company ranging from $[\$ 50,000 \dots \$ 60,000]$ rather than specifying.

$[\$ 51,252.50 \dots \$ 61,252.612]$.

In this we use the method 3-4-5 i.e, the data is partitioned into Equiwidth of 3, 4 or intervals. This method contains the following steps.

Step-1: If the most significant bit is 3, 6, 7 or 9 then the data is partitioned it into 3 equiwidth intervals.

Step-2: If the most significant bit is 2, 4 or 8 then the data is partitioned it into 4 equiwidth intervals.

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Step 3: If the most significant bit is 1, 5 or 10 then the data is partitioned into 5 equiwidth intervals.

For ex, consider the profit at all branches for all electronics company ranging from - \$ 351,986... \$ 4,700,876.50 and also 5% & 95% percentiles are - \$ 159,896 and \$ 1,838,761. It contains the following steps.

Step-1: The given data MIN = - \$ 351,986, MAX = \$ 4,700,876.50. Then the 5% is treated as low and 95% Percentile value is treated as high value

Step-2: If we examine low and the high most significant bit is million dollar bit position i.e. Most significant bit = \$ 1,000,000 (msd) most significant Digit. Rounding low down to significant bit i.e. million dollar bit position

we get low' = - \$ 1,000,000

Similarly, Rounding high upto million dollar bit position we get high' = \$ 2,000,000

$$\Rightarrow (\text{high}' - \text{low}') / \text{msd} = \frac{\$ 2,000,000 - (-1,000,000)}{1,000,000}$$

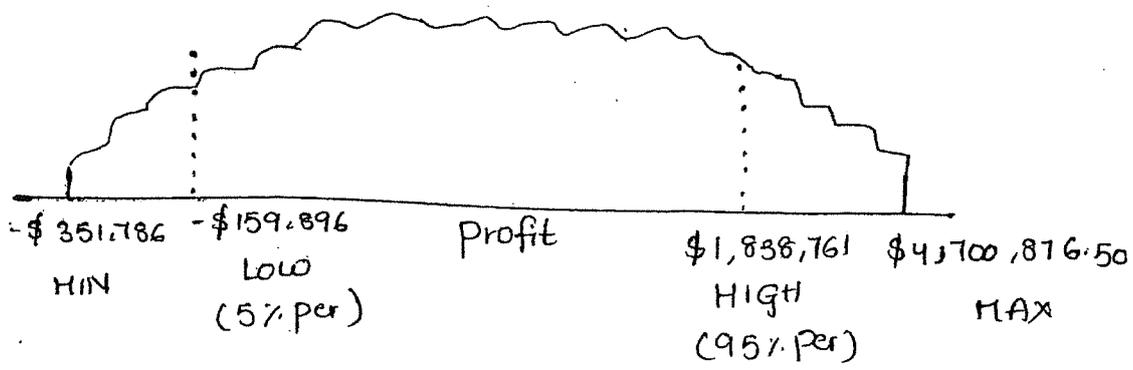
Then we get the new range. This range is partitioned into 3 equiwidths.

Step-3: If we examine MIN and MAX then we get the MIN' = - \$ 4,000,000 and MAX' = \$ 5,000,000.

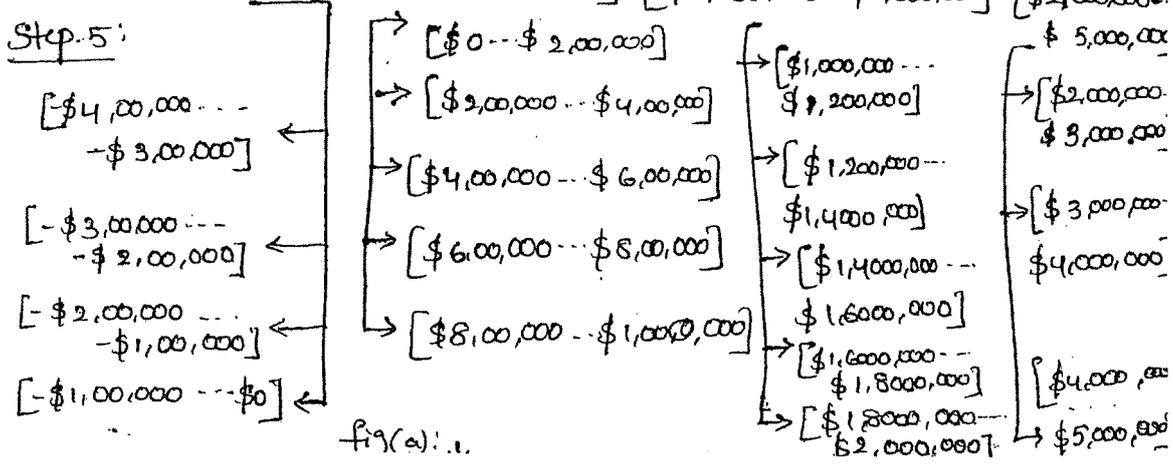
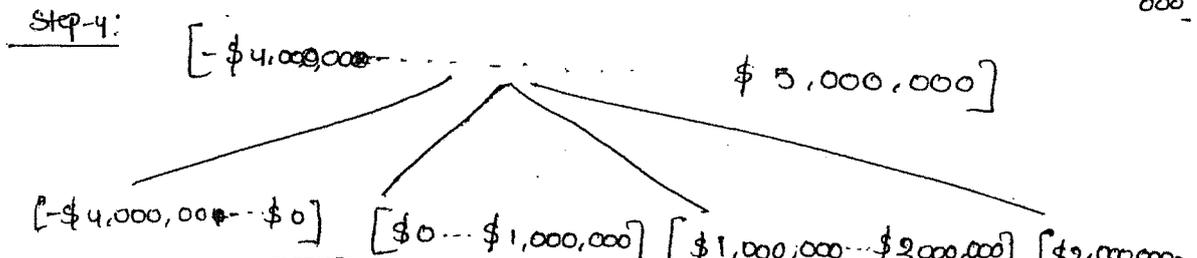
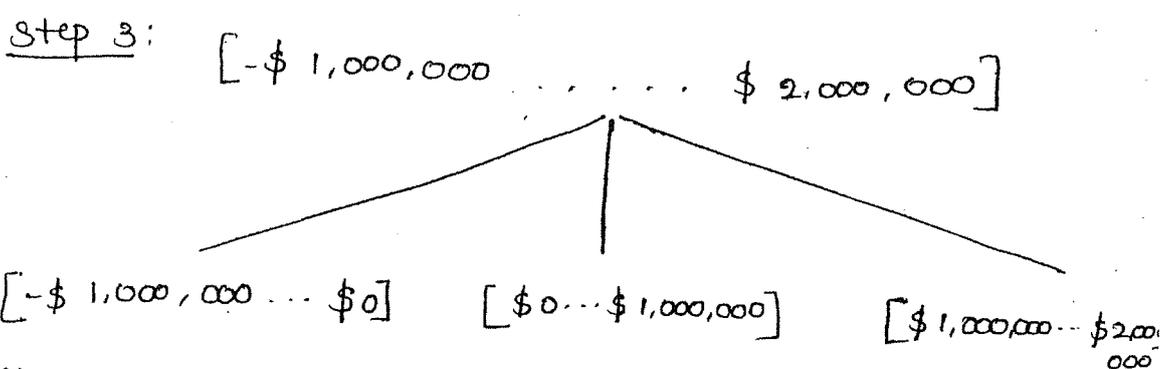
Then we get the new range. This new range is partitioned into 4 equiwidth intervals.

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Step-4: These partitions again partitioned into sub-partitions to define the partition hierarchy. The entire process is shown below.



Step 2: $msd = \$1,000,000$
 $Low' = \$1,000,000$ $HIGH = \$2,000,000$



fig(a)...

Fig(a): Concept hierarchy for Profit based on 3-4-5 rule

Concept Hierarchy Generation for Categorical Data:

The Categorical data means discrete. It contains the finite no. of values for ex, job category, item types etc. It is mainly classified into 2.

1. specification of partial ordering of attributes at the schema level by user:

Here we specify the partial ordering of concept hierarchy at the schema level itself. For ex, Location Dimension contains attributes street, city, State, country. Then the partial.

chapter-4

Getting to know your Data

1.1 Data objects and Attribute types

Data sets are made up of data objects. A data object represents an entity.

For ex^t In Sales DB, the objects may be customers, store items, sales;

medical DB the objects may be patients.

Data objects are typically described by attributes. Data objects can also be referred to as samples, examples, instances, data points or objects.

If the data objects are stored in a database they are data tuples. i.e, the rows of a DB correspond to the data objects, & columns corresponds to the attributes.

What is an Attribute?

An attribute is a data field, representing a characteristic or feature of a data object. The terms attribute, dimension, feature & variable are often used interchangeably in the literature. The term dimension is commonly used in DWH. In Machine learning, we use the term feature, while statisticians refer the term variable.

Observed values for a given attribute are known as observations. A set of attributes used to describe a given object is called an attribute vector.

Types of attributes:-

The type of an attribute is determined by the set of possible values. they are

- 1) nominal
- 2) binary
- 3) ordinal
- 4) numeric.

1) Nominal Attributes

Nominal means "relating to names". The values of a nominal attribute are symbols or names of things. Each value represents some kind of category, code (or) state. Nominal attributes are also referred to as categorical. The values do not have any meaningful order. In computer science, the values are known as enumerations.

Ex 1 -> Suppose that hair_color & marital_status are

- two attributes describing person objects.

Possible values for hair_color are black, brown, red, gray, white.

The attribute marital-status can take on the values single, married, divorced & widowed.

2, Binary Attributes —

Binary attr. is a nominal attr. with only two categories or states: 0 (or 1), where 0 represents absent of an attribute, & 1 represents present of an attr. Binary attr. are referred to as Boolean if the two states correspond to true & false.

for ex: Given the attribute Smoker describing a patient object, 1 indicates that the patient smokes, while 0 indicates that the patient does not.

A binary attribute is symmetric if both of its states are equally valuable and costly & costly the same weight; i.e., there is no preference on which outcome should be coded as 0 or 1.

A binary attribute is Asymmetric if the outcomes of the states are not equally important. Such as the positive & negative outcomes of a medical test for HIV. By convention, we code the most important outcome, which is usually the rarest one, by 1 & the other by 0.

3, ordinal Attributes :-

An ordinal att_i is an att_i with possible values that have a meaningful order or ranking among them, but the magnitude b/w successive values is not known.

~~Ex~~ Ordinal att_s are useful for registering subjective assessments of qualities that cannot be measured objectively. Ordinal attributes are often used in surveys for rating.

Ex :- In one survey, participants were asked to rate how satisfied they were as customer. Customer satisfaction had the following ordinal categories.

- 0: very dissatisfied 1: somewhat dissatisfied
- 2: neutral 3: satisfied & 4: very satisfied.

4, Numeric Attributes :-

A numeric attribute is quantitative; i.e., it is a measurable quantity, represented in integer or real values. Numeric att_s can be interval-scaled or ratio-scaled.

Interval-scaled Attributes

Interval-scaled att_s are measured on a scale of equal-size units. The values of interval-scaled

attributes have order and can be positive, 0, or -ve.

In addition to providing a ranking of values, such attributes allow us to compare & quantify the different values.

ex:

- 1) Celsius temperature
- 2) IQ (Intelligence scale)
- 3) Time on a clock with hands.

Ratio-Scaled Attributes

A ratio-scaled attribute is a numeric attribute with an inherent zero point. i.e., if a measurement is ratio scaled, we can speak of a value as being a multiple of another value. In addition, we can also compute the different values, as well as mean, median & mode.

ex:

- 1) Age, weight, Height.
- 2) Ruler measurements
- 3) Years of education.

5) Discrete versus Continuous Attributes :-

In we have organized attr: into nominal, binary, ordinal, & numeric types. there are many ways to organize attr: types. the types are not mutually exclusive.

Discrete attribute :-

It has a finite or countably infinite set of values, which may or may not be represented as integers. The attr: hair-color, smoker, medical-test each have a finite number of values, such as ~~one~~ for binary attr: and are discrete.

Continuous Attribute :-

If an attr: is not discrete, then it is continuous. The terms numeric attribute & continuous attr: are often used interchangeably in literature. Real values are represented using a finite number of digits. Continuous attributes are typically represented as floating point variables.

9.2. Basic Statistical Description :-

Basic statistical descriptions can be used to identify properties of the data and highlight which data values should be treated as noise or outliers.

In this we discuss three areas of basic statistical descriptions. measures of central tendency, measures of dispersion & graphic displays.

9.2.1 Measuring the Central Tendency : mean, median and Mode:

A measure of central tendency is a single value that describes the way in which a group of data cluster around a central value. In other words, it is a way to describe the centre of a data set. There are three measures of central tendency:

- 1) Mean
- 2) Median
- 3) Mode.

Mean

The mean is preferred measure of central tendency because it considers all the values of dataset. In order to calculate the mean, data must be numerical. You cannot use the mean for nominal data, which is data on characteristics like gender.

The mean for set of values is

$$\bar{x} = \frac{\sum_{i=1}^N x_i}{N} = \frac{x_1 + x_2 + \dots + x_N}{N}$$

Ex Suppose we have the following values for salary, shown in increasing order:

30, 36, 47, 50, 52, 52, 56, 60, 63, 70, 70, 110

$$\bar{x} = \frac{30 + 36 + 47 + 50 + 52 + 52 + 56 + 60 + 63 + 70 + 70 + 110}{12}$$

$$= \frac{696}{12}$$

$$= 58$$

The mean salary is \$58,000.

a)

Sometimes, each value x_i in a set may be associated with a weight w_i for $i = 1, 2, \dots, N$. The weights reflect the significance.

$$\bar{x} = \frac{\sum_{i=1}^N w_i x_i}{\sum_{i=1}^N w_i} = \frac{w_1 x_1 + w_2 x_2 + \dots + w_N x_N}{w_1 + w_2 + \dots + w_N}$$

This is called as weighted average or weighted Arithmetic mean (AM).

Median:-

The median is the middle value for the given data, that has been arranged in order of magnitude.

If we have 2 middle values, we can find the mean value b/w the 2 values, & take it as a median.

The median is expensive to compute when we have a large number of observations.

$$\text{Median} = l + \left(\frac{\frac{n}{2} - cf}{f} \right) \times h$$

l = lowest values (lowest interval)

$$\frac{n}{2} = \frac{\text{no. of observations}}{2}$$

cf = highest frequency

cf = above the f

h = class interval.

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Mode

A statistical term that refers to the most frequently occurring number found in a set of numbers. The mode is found by collecting and organizing the data in order to count the frequency of each result.

Ex:

Suppose we have the following values

30, 36, 47, 50, 52, 52, 56, 60, 63, 70, 70, 70, 110

mode = 70, 52. This is bimodal.

For unimodal numeric data that are moderately skewed, we have the following expression.

$$\text{mean} - \text{mode} \approx 3 \times (\text{mean} - \text{median})$$

i.e. the mode for unimodal frequency curves that are moderately skewed can easily be approximated if the mean & median values are known.

Mid range:
Range

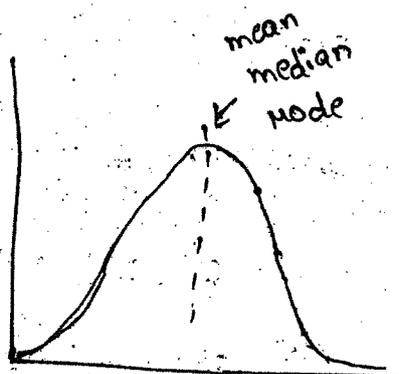
(the range of a set of data is the diff^s b/w largest & smallest values)

For the above example the midrange is as follows

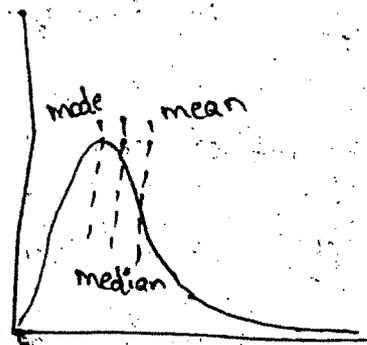
$$\frac{30,000 + 110,000}{2} = 70,000$$

In a unimodal frequency curve with perfect symmetric data distribution the mean, median & mode are all at the same center value.

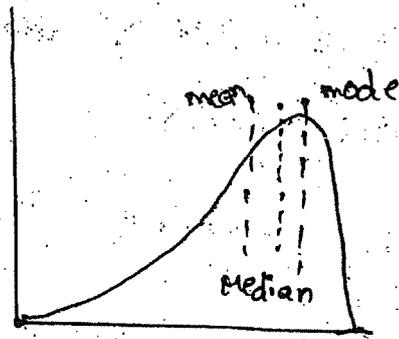
Data in most real application are not symmetric. They may instead be either "positively skewed", where the mode occurs at a value ~~th~~ smaller than median, or negatively skewed, where the mode occurs at a value greater than the median.



(a) Symmetric data



(b) Positively skewed data



(c) Negatively skewed data.

fig (a.1) Mean, median, mode of symmetric v/s positively & negatively skewed data.

Q.2 Measuring the Dispersion of Data :-

A measure of spread, or measure of dispersion is used to describe the variability in a small sample or population. The measures include range, quantiles, quartiles, percentiles and the interquartile range. The five number summary, which can be displayed as a boxplot, is useful to identifying outliers.

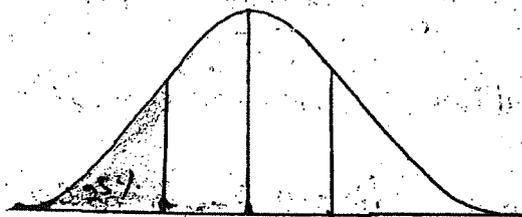
Range, quartiles, and Interquartiles range

Let x_1, x_2, \dots, x_N be a set of observations for some numeric attribute, X . The range of the set is the difference b/w the large & small values.

$$\text{range} = \max(x) - \min(x).$$

Quartiles tell us about the spread of the data set by breaking the data set into quarters, just like median breaks it in half.

The first quartile (Q_1) lies b/w the 25th & 26th observation, the second quartile (Q_2) b/w 50th & 51st observation & the third quartile (Q_3) b/w the 75th & 76th observation.



Q_1 Q_2 Q_3
25th median 75th
percentile percentile.

or

The distance b/w the first & third quartiles is simple measure of spread that gives the range covered by the middle half of the data.

This distance is called "Interquartile range" (IQR)

$$IQR = Q_3 - Q_1$$

Ex 21 Suppose we have have the following values for salary is

30, 36, 47, 50, 52, 52, 56, 60, 63, 70, 70, 110.

It contains 12 observations, already stored in increasing order. Thus, the quartiles for the data are 3rd, 6th & 9th values respectively.

$$Q_1 = 47,000$$

$$Q_2 = 52,000$$

$$Q_3 = 63,000$$

$$IQR = 63 - 47 = 16,000.$$

five-number Summary, Boxplots and outliers:-

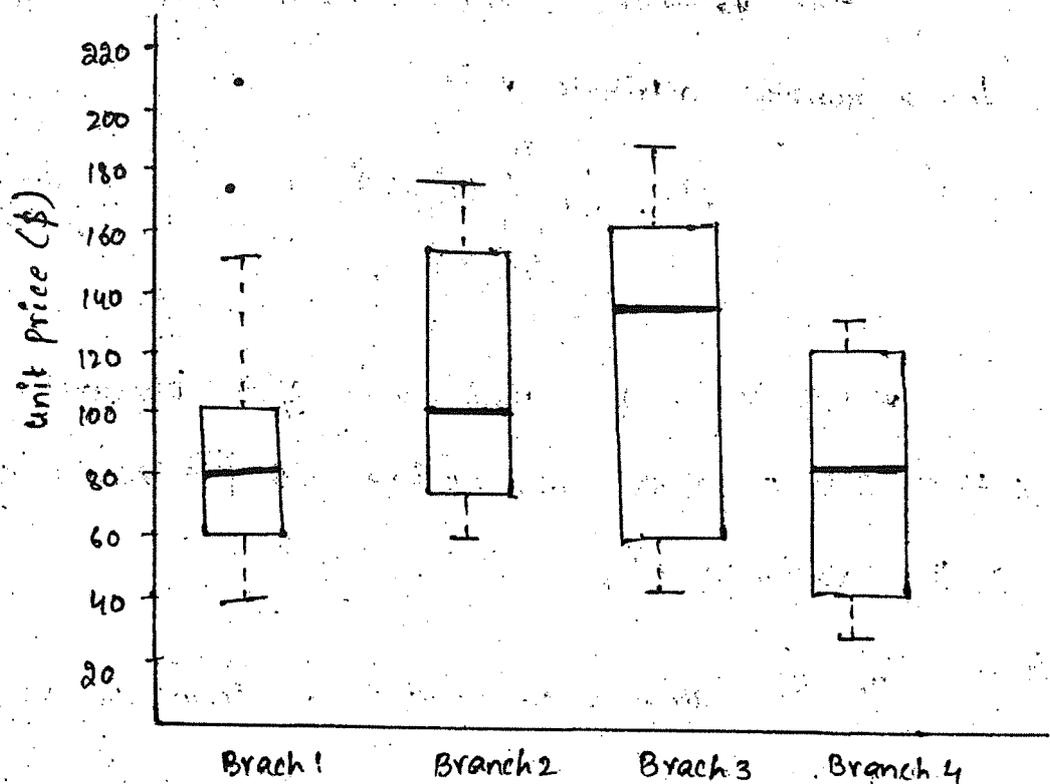
The outlier is an observation that lies an abnormal distance from other values.

A common rule of thumb for identifying suspected outliers is to single value falling at least $1.5 \times IQR$ above the 3rd quartile or below the 1st quartile.

The five-number summary of a distribution consists of the median (Q_2), the quartiles Q_1 & Q_3 , and the smallest & largest individual observations written in the order of min, Q_1 , median, Q_3 , max.

Box plots are a popular way of visualizing a distribution. A boxplot incorporates the five number summary as follows:

- * The ends of the box are the the quartiles, so the box length is interquartile range.
- * The median is marked by a line within the box.
- * Two lines (called whiskers) outside the box extend to the smallest & largest observations.



Q.3 Box plot for the unit price data for items sold at four branches of all electronics during a given time period.

Variance and Standard Deviation:

Variance and standard deviation are measures of data dispersion. They indicate how spread out a data distribution is.

A low standard deviation means that the data observations tend to be very close to the mean, while high S.D indicates that the data are spread out over a large range of values.

The variance of N observations x_1, x_2, \dots, x_N for a numeric attribute x is

$$\begin{aligned}\sigma^2 &= \frac{1}{N} \sum_{i=1}^N (x_i - \bar{x})^2 \\ &= \left(\frac{1}{N} \sum_{i=1}^N x_i^2 \right) - \bar{x}^2\end{aligned}$$

where \bar{x} is the mean value of the observations.

If the S.D σ , of the observations is the square root of the variance σ^2 .

Ex 1 In the above observation we found $\bar{x} = 58,000$ and $N = 12$.

$$\begin{aligned}\sigma^2 &= \frac{1}{12} (30^2 + 36^2 + \dots + 110^2) - 58^2 \\ &= 379.17\end{aligned}$$

$$\sigma = \sqrt{379.17}$$

$$\approx 19.47$$

The basic properties of S.D σ , as a measure of spread are as follows:

- 1) σ measures spread about the mean and should be considered only when the mean is chosen as the measure of center.
- 2) $\sigma = 0$ only, when there is no spread, i.e., when all observations have same value. otherwise $\sigma > 0$.

4.2.3 Graphic Displays of Basic Statistical Descriptions of Data :-

The graphic displays of basic statistical descriptions include quantile plots, quantile-quantile plots, histograms & scatter plots. Such graphs are useful for the visual inspection of data, which is useful for data preprocessing. First three of these show univariate distributions, and while scatter plots show bivariate distributions.

Quantile plot:-

A quantile plot is a simple, well known and effective way to have a first look at a univariate

data distribution. It displays all of the data for the given attribute. then, it plots quantile information.

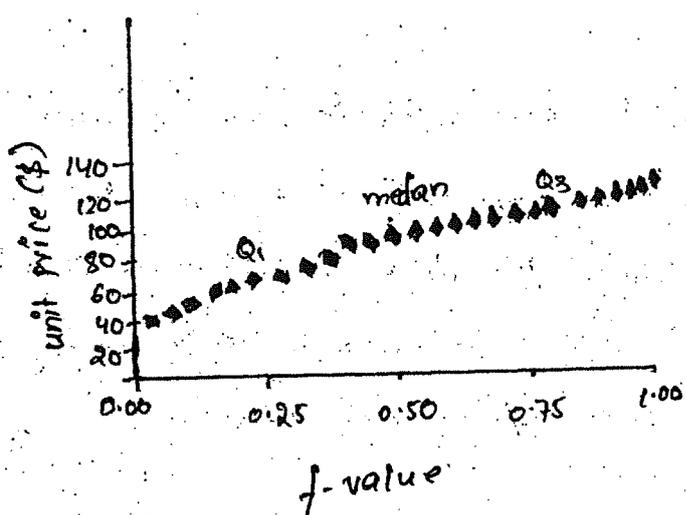
ex: let x_i , for $i=1$ to N , be the ^{data} sorted in increasing order so that x_1 is the smallest element of x and x_N is the largest for some ordinal or numeric attribute x .

A set of unit price data for items sold at a Branch of All Electronics

unit price (\$)	count of items sold
40	275
43	300
47	250
-	-
74	-
75	360
78	515
-	540
115	-
119	320
120	270
	350

Table A.1

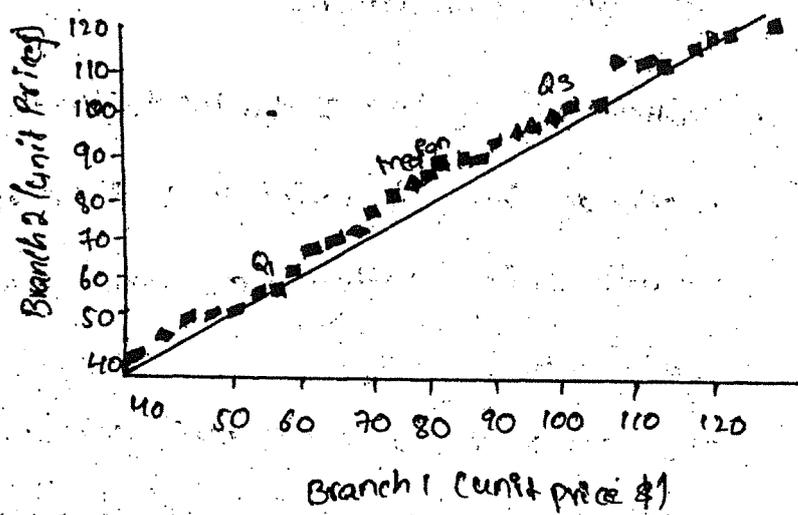
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A quantile plot for the unit price data of table 21.

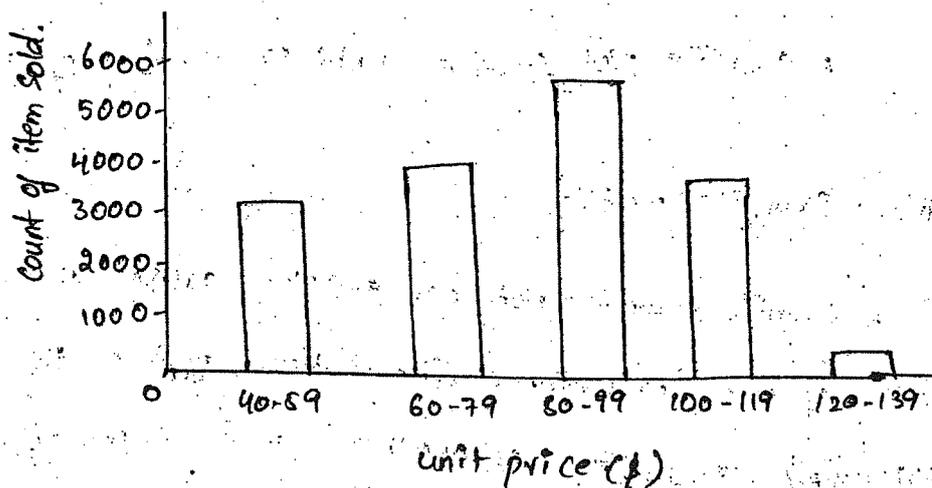
Quantile - Quantile plot:-

A quantile - quantile plot, or q-q plot, graphs the quantiles of one univariate distribution against the corresponding quantiles of another. It is a powerful visualization tool in that it allows the user to view whether there is a shift in going from one distribution to another.



Histograms

A histogram is an accurate representation of numerical data. To construct a histogram for an attribute x , partitions the data distributions of x into buckets. Each bucket is represents only a single attribute of value / frequency pair.

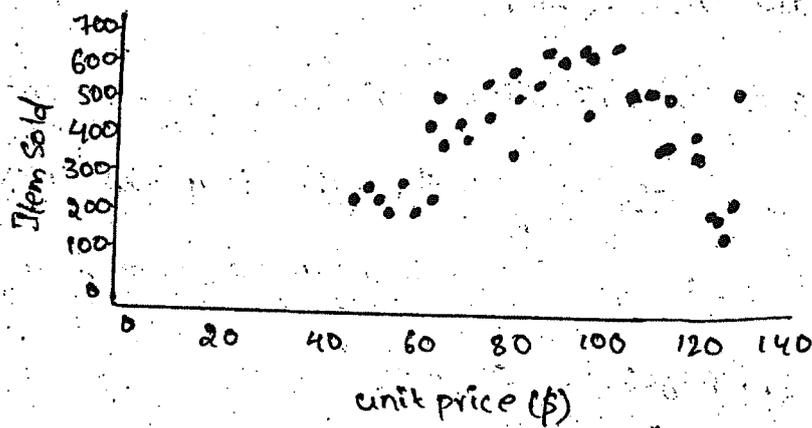


A histogram for table 4.1

Scatter plots and Data correlation

A scatter plot is one of the most effective graphical methods for determining if there appears to be a relationship, pattern or trend b/w two numeric attributes.

To construct a scatter plot, each pair of values is treated as a pair of coordinates in an algebraic and plotted as points in the plane.



scatter plot for the table 2.1

The scatter plot is useful to represent the bivariate data to clusters of points & outliers, or to explore the possibility of correlation relations.

Two attributes, X , and Y , are correlated if one attribute implies to other. Correlation can be +ve, -ve or null.



(a) positive



(b) negative

Scatter plots can be used to find (a) positive or (b) negative b/w attributes.

values

c

* 4.3 Data Visualization

Data visualization aims to communicate data clearly and effectively through graphical represents.

Data visualization has been used extensively in many applications of

for ex: at work for reporting, managing business operations, and tracking progress of tasks.

more popularly, we can take advantage of visualization techniques to discover data relationships that are otherwise not easily observable by looking at the raw data.

4.3.1 Pixel oriented visualization techniques:-

A simple way to visualize the value of dimension is use a pixel where the colors of the pixel reflect the dimension's value. For the data set of m dimensions, pixel-oriented technique creates m windows on the screen, one for each dimension.

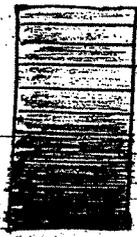
ex: All Electronics maintains a customer info table which consists of income, credit limit, transaction

Volume and age. we can analyze correlation b/w income and the other attributes by visualization?

we can sort all the customers in income ascending order, & used this order to layout the customer data in four visualization windows, as follows.



(a) income



(b) credit limit



(c) transaction - volume



(d) age

Pixel oriented visualization of four attr.

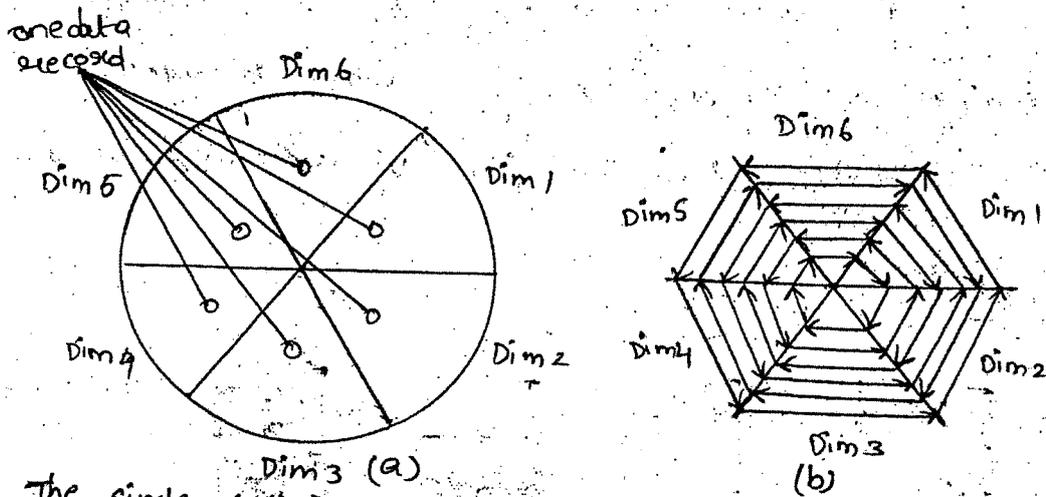
Credit limit increases as income increases; whose income is in the middle range are more likely to purchase more from All Electronics; there is no clear correlation b/w income & age.

Filling a window by laying out the data records in a linear way may not work well for a wide window. we can lay out the data records

in a space-filling curve to fill the windows.

A SF is a curve with a range that covers the entire n-dimensional unit hypercube. Since the visualization

windows are 2-D, we can use any 2-D space filling curve as follows.



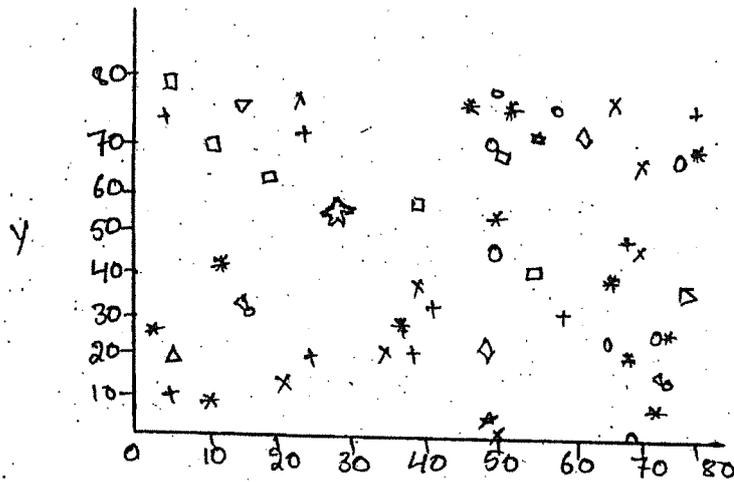
The circle segment technique. (a) Representing a data record in circle segments (b) laying out pixels in circle segments.

4.3.2 Geometric projection visualization techniques :-

Geometric projection techniques help users find interesting projections of multidimensional data sets. The central challenge the geometric projection techniques try to address is how to visualize a high dimen- sional space on a 2-D display.

A scatter plot displays 2-D data points using Cartesian coordinates. A third dimension can be added using different colors or shapes to shapes to represent diff. data points.

11/19



Visualization of a 2-D data set using a scatter plot.

where x and y are two spatial attributes and the third dim₃ is represented by different slopes.

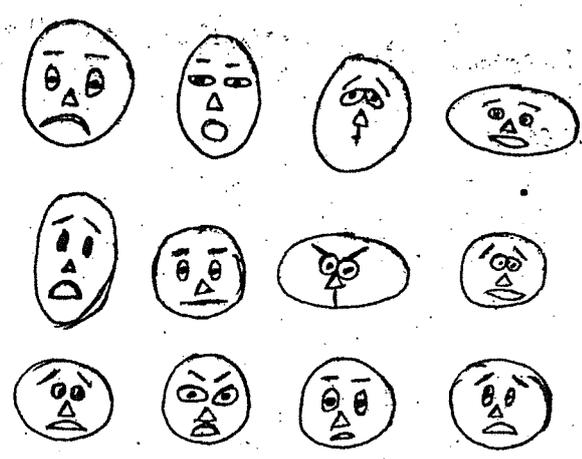
Through this visualization, we can see that the points of type "+" and "x".

A scatter plot matrix technique is a useful extension to the scatter plot. For an n -dimensional data set, a scatter plot matrix is an $n \times n$ grid of 2-D scatter plots that provides a visualization of each dim_i with every other dim_j. The scatter plot matrix becomes less effective as the dimensionality increases.

4.3.3 :- Icon Based Visualization Techniques

Icon-based visualization techniques use small icons to represent multidimensional data values. There are two popular icon-based techniques: ^{chernoff} chernoff faces and stick figures.

Chernoff faces were introduced in 1973 by statistician Herman Chernoff. They display multidimensional data of up to 18 variables as a cartoon human face.



Chernoff faces. Each face represents an n-dimensional

data point ($n \leq 18$)

4.3.4 Hierarchical Visualization Techniques :-

The visualization techniques discussed so far focus on visualizing multiple dimensions simultaneously. However, for a large data set of high dimensionality, it would be difficult to visualize all dimensions at the same time.

Hierarchical visualization techniques partition all dimensions into subsets. The subsets are visualized in a hierarchical manner.

~~works~~ "worlds-within-worlds" also known as n-vision, is a representative hierarchical visualization method. Suppose we want to visualize a 6-D dataset, where the dimensions are F, X_1, \dots, X_5 . We can first fix the values of dimensions X_3, X_4, X_5 to some selected values, say c_3, c_4, c_5 . We can then visualize F, X_1, X_2 using 3-D plot.

4.3.5 Visualizing Complex Data and Relations :-

Visualization techniques were mainly for numeric data. Recently, more and more non-numeric data, such as text and social networks, have become available.

For ex, many people on the web tag various objects such as pictures, blog entries and product reviews.

4.4 Measuring Data Similarity and Dissimilarity :-

In data mining applications, such as clustering, outliers analysis, and nearest-neighbour classification, we need ways to assess how ~~like~~ alike or unlike objects are in comparison to one another.

For ex, a store may want to search for clusters of customer objects, resulting in group of customers with similar characteristics. (income, age of residence page). Such information can then be used for marketing.

A cluster is a collection of data objects such a cluster are similar to one the objects.

so employs clustering-based potential outliers as objects to others.

ct similarities can also be used ation schemes where a given s label based on its similarity the model.

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M.C.A. DEGREE EXAMINATIONS
FOURTH SEMESTER

Paper - III : DATA MINING CONCEPTS AND
TECHNIQUES

(W.E.F. 2019-20 Admitted Batch)
Time : 3 Hours

Maximum : 75 Marks

SECTION - A

1. Answer all questions.

a. What is multi -

dimensional analysis? How it is implemented through OLAP? Explain various types of OLAP systems used for multi - dimensional data.

(4×15=60)

4.4.1 Data Matrix versus Dissimilarity Matrix :-

We studying the central tendency, dispersion & spread of observed values for some attribute x . that are one-dimensional, i.e, described by a single attribute. we talk about objects described by multiple attributes. Therefore, we need a change in notation.

suppose that we have n objects (ex: persons, items, or courses) described by p attributes (age, height, weight, or gender). The objects are $x_1 = (x_{11}, x_{12}, \dots, x_{1p})$, $x_2 = (x_{21}, x_{22}, \dots, x_{2p})$ and so on, where x_{ij} is the value for object x_i of the j^{th} attribute.

Main memory-based clustering and nearest-neighbour algorithms typically operate on either of the following two data structures.

1) Data matrix (object-by-attribute structure) :-

this is also called as object-by-attribute structure. this structure stores the n data objects in form of relational table, or $n \times p$ matrix (n objects \times p attributes).

x_{11}	\dots	x_{1f}	\dots	x_{1p}
\dots	\dots	\dots	\dots	\dots
x_{i1}	\dots	x_{if}	\dots	x_{ip}
\dots	\dots	\dots	\dots	\dots
x_{n1}	\dots	x_{nf}	\dots	x_{np}

2) Dissimilarity matrix :-

This is also called as object-by-object structure. This structure stores a collection of proximities that are available for all pairs of objects. It is also represented as $n \times n$ table.

$$\begin{bmatrix} 0 & & & & \\ d(2,1) & 0 & & & \\ d(3,1) & d(3,2) & 0 & & \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ d(n,1) & d(n,2) & \dots & \dots & 0 \end{bmatrix}$$

where $d(i,j)$ is the measured dissimilarity (or) difference b/w objects i & j . $d(i,j)$ is a non-ve number that is close to 0 when objects i & j are highly similar or near each other.

$d(i,i) = 0$ i.e., the difference b/w an object and itself is 0.

Measure of similarity can be expressed as a function of measures of dissimilarity.

Ex :- for nominal data

$$\text{sim}(i,j) = 1 - d(i,j)$$

where $\text{sim}(i,j)$ is the similarity b/w objects i and j .

A data matrix is made up of two entities they are rows (for objects) and columns (for attributes). Therefore, the data matrix is often called as "two-mode" matrix. The dissimilarity matrix contains one kind of entity & so is called as "one-mode" matrix.

4.4.2 Proximity measures for Nominal Attributes:-

A nominal attribute can take on two or more states. For example: map-color is a nominal attribute that may have say five states: red, yellow, green, pink and blue.

Let the no. of states of a nominal attribute be M . The states can be denoted by letters, symbols or a set of integers, such as $1, 2, \dots, M$.

dissimilarity b/w objects described by nominal attributes

The dissimilarity b/w two objects i and j can be computed based on the ratio of mismatches:

$$d(i, j) = \frac{p - m}{p}$$

where m is the no. of matches i.e. no. of attributes for which i & j are in the same state. and p is the total no. of attributes describing the objects.

Ex: Suppose that we have the sample data of as follows, expect that only the object-identifier and the attribute test-1 are available, where test-1 is nominal.

Table 4.2. A simple Data Table Containing Attributes of mixed type.

Object Identifier	test-1 (nominal)	test-2 (ordinal)	test-3 (numeric)
1	code A	excellent	45
2	code B	fair	22
3	code C	good	64
4	code A	excellent	28

Let's compute the dissimilarity matrix.

$$\begin{bmatrix} 0 & & & \\ d(2,1) & 0 & & \\ d(3,1) & d(3,2) & 0 & \\ d(4,1) & d(4,2) & d(4,3) & 0 \end{bmatrix}$$

we set $p=1$, so that $d(i,j)$ evaluates to 0 if the i 's are match, & $\neq 1$ if the objects differ. Thus we get

$$\begin{bmatrix} 0 & & & \\ 1 & 0 & & \\ 1 & 1 & 0 & \\ 0 & 1 & 1 & 0 \end{bmatrix}$$

we see all the objects are dissimilar except objects 1 and 4 i.e. $d(4,1) = 0$.

Alternatively, similarity can be computed as

$$\text{sim}(i, j) = 1 - d(i, j) = \frac{m}{p}$$

4.4.8:- Proximity Measure for Binary Attributes:-

A binary attribute has only one of two states: 0 or 1 where 0 means that the attribute is absent, and 1 means that it is present. Treating binary attributes as if they are numeric can be misleading. Therefore, methods specific to binary data are necessary for computing dissimilarity.

One approach involves computing a dissimilarity matrix from the given binary data. If all binary attributes are thought of as having the same weight, we have the 2x2 contingency table as follows

		object j		sum
		1	0	
object i	1	q	r	q+r
	0	s	t	s+t
sum		q+s	r+t	p

Contingency Table for Binary Attributes.

where 'q' is the number of attributes that equal to 1 for both objects i & j, 'r' is the no. of attributes that equal '1' for object i, but equal '0' for object j, 's' is the number of attributes that equal '1' for object 'j' but equal 0 for object i, and 't' is the no. of attributes that equal '0' for both objects i & j. The total no. of attributes is P

$$P = q + r + s + t.$$

For Symmetric binary attributes, each state is equally valuable. Dissimilarity that is based on symmetric binary attributes is called symmetric binary dissimilarity. If objects i and j are described by symmetric binary attributes, then the dissimilarity

between i & j is

$$d(i, j) = \frac{r + s}{q + r + s + t}$$

For asymmetric binary attributes, the two states are not equally important, given two asymmetric binary attributes, the agreement of two 1's is then considered more significant than that of two 0's. Therefore, such binary attributes are often considered "monary" (having one state). The dissimilarity based on these attributes is called asymmetric binary dissimilarity.

$$d(i,j) = \frac{r+s}{2+r+s}$$

where the no. of negative matches t , is considered unimportant and is thus ignored.

when both Symmetric & asymmetric binary attributes occur in the same data set, the mixed attributes approach can be applied.

Ex 1

Suppose the patient record table contains the attribute name, gender, fever, cough, test-1, test-2, test-3, & test-4.

where name is an object identifier, gender is a symmetric attribute, and the remaining attributes are asymmetric binary.

name	gender	fever	cough	test-1	test-2	test-3	test-4
Jack	M	Y	N	P	N	N	N
Jim	M	Y	Y	N	N	N	N
Mary	F	Y	N	P	N	P	N
⋮	⋮	⋮	⋮	⋮	⋮	⋮	⋮

4.4 Relational Table where patients are described by Binary Attributes.

For Asymmetric attribute values, the values q and p be set to 1, and the value n (no or negative) be set to '0'.

The distance b/w each pair of the three patients Jack, Mary & Jim is

$$d(\text{Jack}, \text{Jim}) = \frac{1+1}{1+1+1} = 0.67$$

$$d(\text{Jack}, \text{Mary}) = \frac{0+1}{2+0+1} = 0.33$$

$$d(\text{Jim}, \text{Mary}) = \frac{1+2}{1+1+2} = 0.75$$

These measurements suggest that Jim & Mary are unlikely to have a similar disease. and Jack and Mary are the most likely to have a similar disease.

4.4.4 Dissimilarity of Numeric Data: Minkowski Distance

We describe distance measures that are commonly use for computing the dissimilarity of objects by numeric attributes. These measures include the Euclidean, Manhattan and Minkowski distances.

In some cases, the data are normalized before applying distance calculations. This involves transforming the data to fall within a smaller or common range such as $[-1, 1]$ or $[0, 1]$.

3)
 4)
 The most popular distance measure is "Euclidean distance". Let $i = (x_{i1}, x_{i2}, \dots, x_{ip})$ and $j = (x_{j1}, x_{j2}, \dots, x_{jp})$ be two objects described by p numeric attributes. The Euclidean distance b/w objects i & j is

$$d(i, j) = \sqrt{(x_{i1} - x_{j1})^2 + (x_{i2} - x_{j2})^2 + \dots + (x_{ip} - x_{jp})^2}$$

ex: Let $x_1 = (1, 2)$ & $x_2 = (3, 5)$, the Euclidean distance b/w two objects is

$$d(x_1, x_2) = \sqrt{(1-3)^2 + (2-5)^2} = \sqrt{2^2 + 3^2} = 3.61$$

Another well-known measure is the Manhattan distance or city block distance, name so because it is the distance in blocks b/w any two points in a city.

$$d(i, j) = |x_{i1} - x_{j1}| + |x_{i2} - x_{j2}| + \dots + |x_{ip} - x_{jp}|$$

ex: Manhattan distance for the above example

$$d(x_1, x_2) = |1-3| + |2-5| = 2+3 = 5$$

Both the Euclidean and the Manhattan distance satisfy the following mathematical properties.

- 1) Non-negativity: $d(i, j) \geq 0$. : Distance is a non-ve number.
- 2) Identity of indiscernibles: $d(i, i) = 0$. The distance of an object to itself is '0'.

3) Symmetry: $d(i, j) = d(j, i)$: Distance is a symmetric function.

4) Triangle inequality: $d(i, j) \leq d(i, k) + d(k, j)$.

"Minkowski distance" is a generalization of the Euclidean and Manhattan distances.

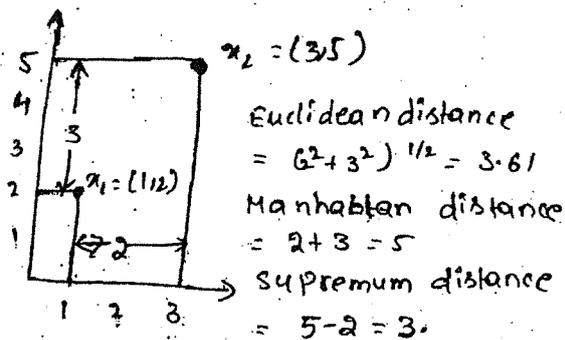
$$d(i, j) = \sqrt[h]{|x_{i1} - x_{j1}|^h + |x_{i2} - x_{j2}|^h + \dots + |x_{ip} - x_{jp}|^h}$$

where h is a real number such that $h \geq 1$. It represents the Manhattan distance when $h=1$ and Euclidean distance when $h=2$.

The supremum distance is a generalization of the Minkowski distance for $h \rightarrow \infty$. To compute it, we find the $\text{arg } f$ that gives the maximum difference in values b/w two objects.

$$d(i, j) = \lim_{h \rightarrow \infty} \left(\sum_{f=1}^p |x_{if} - x_{jf}|^h \right)^{1/h}$$

$$= \max_f |x_{if} - x_{jf}|$$



Euclidean, Manhattan, supremum distance between two objects

4.4.5 Proximity Measure for Ordinal Attributes :-

4.4

The values of an ordinal attributes have a meaningful order or ranking. The treatment of ordinal attributes is quite similar to that of numeric attributes when computing dissimilarity b/w two objects.

Suppose f is an attribute from the set of ordinal attributes describing n objects. The dissimilarity computation w.r. to f involves the following steps:

1. The value of ' f ' for i th object is x_{if} , and f has M_f ordered states, representing the ranking $\{1, \dots, M_f\}$. Replace each x_{if} by its corresponding rank $r_{if} \in \{1, \dots, M_f\}$.
2. We perform data normalization by replacing the rank r_{if} of the i th object in the f th attribute by
$$z_{if} = \frac{r_{if} - 1}{M_f - 1}$$
3. Dissimilarity can be computed using any of the distance measures described in numeric attributes, using z_{if} to represent the f value for i th object.

*

4.4.6 Dissimilarity for Attributes of Mixed Types :-

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es

One approach is to group each type of attribute together, performing separate data mining analysis for each type eg: clustering. This is feasible if these analyses derive compatible results.

ly

Suppose that the data set contains p attr's of mixed type. The dissimilarity b/w i & j objects is defined as

$$d(i,j) = \frac{\sum_{f=1}^p \delta_{ij}^{(f)} \cdot d_{ij}^{(f)}}{\sum_{f=1}^p \delta_{ij}^{(f)}}$$

3

where the indicator $\delta_{ij}^{(f)} = 0$ if either

- 1) x_{if} or x_{jf} is missing. (or)
- 2) $x_{if} = x_{jf} = 0$ and attribute f is asymmetric binary.

otherwise, $\delta_{ij}^{(f)} = 1$.

The contribution of attribute f to the dissimilarity b/w i & j i.e., $d_{ij}^{(f)}$ is computed dependent on its type:

* If f is numeric $d_{ij}^{(f)} = \frac{|x_{if} - x_{jf}|}{\max_h x_{hf} - \min_h x_{hf}}$

where h runs over all nonmissing objects for attr f .

* If f is nominal or binary: $d_{ij}^{(f)} = 0$ if $x_{if} = x_{jf}$;
otherwise $d_{ij}^{(f)} = 1$.

* If f is ordinal: compute the ranks r_{ij} and

$$z_{ij} = \frac{r_{ij} - 1}{M_j - 1}, \text{ and treat } z_{ij} \text{ as numeric.}$$

4.4.7 Cosine Similarity :-

A document can be represented by thousands of attributes, each recording the frequency of a particular word or phrase in the document. Thus, each document is an object represented by a "term-frequency vector".

Cosine Similarity is a measure of similarity that can be used to compare documents or give a ranking of documents w.r. to a given vector of query words.

Let x & y be two vectors for comparison. Using cosine measure as a similarity function, we have

$$\text{sim}(x, y) = \frac{x \cdot y}{\|x\| \|y\|}$$

where $\|x\|$ is the Euclidean norm of vector

$x = (x_1, x_2, \dots, x_p)$ defined as $\sqrt{x_1^2 + x_2^2 + \dots + x_p^2}$. $\|y\|$

$\|y\|$ is the Euclidean norm of vector y .

Ex:

— suppose that x & y are the first two

term-frequency vectors in the below table. i.e

$$x = (5, 0, 3, 0, 2, 0, 0, 2, 0, 0) \text{ \& } y = (3, 0, 2, 0, 1, 1, 0, 1, 0, 1)$$

Document Vector or Term-Frequency Vector

Document	team	coach	hockey	baseball	soccer	penalty	score
Document 1	5	0	3	0	2	0	0
Document 2	3	0	2	0	1	1	0
Document 3	0	7	0	2	1	0	0
Document 4	0	1	0	0	1	2	2

win	loss	season
2	0	0
1	0	1
3	0	0
0	3	0

compute the cosine similarity b/w 2 vectors x & y

$$x \cdot y = 5 \times 3 + 0 \times 0 + 0 \times 3 + 2 \times 0 + 2 \times 1 + 0 \times 1 + 0 \times 0 + 2 \times 1 + 0 \times 0 + 0 \times 1 = 25$$

$$\|x\| = \sqrt{5^2 + 0^2 + 3^2 + 0^2 + 2^2 + 0^2 + 0^2 + 2^2 + 0^2 + 0^2} = 6.48$$

$$\|y\| = \sqrt{3^2 + 0^2 + 2^2 + 0^2 + 1^2 + 1^2 + 0^2 + 1^2 + 0^2 + 1^2} = 4.12$$

$$\text{sim}(x, y) = \frac{25}{6.48 \times 4.12} = 0.94$$

Therefore, if we were using the cosine similarity measure to compare these documents, they would be considered quite similar.



5th unit Architecture of Data Mining System.

To get the efficient architecture of DMS. The DMS must be integrated with db (or) DWH. If the DMS is integrated with db or DWH then, we have to consider the following coupling schemas.

- ① No coupling
- ② loosely coupling
- ③ Semitight coupling
- ④ Tight coupling.

① No coupling:- Here, DMS does not integrate with db or DWH. These types of systems contain the following drawbacks.

① If the DMS doesn't integrate with the db then it doesn't provide any performance, scalability etc...

② If the DMS doesn't integrate with the DWH then it doesn't provide data cleaning, data transformation, data integration etc...

These systems extract the data from the files because these systems

doesn't contain any data structures or any algorithms & the results are send to file.

∴ No coupling means poor design.

2. Loosely Coupling:

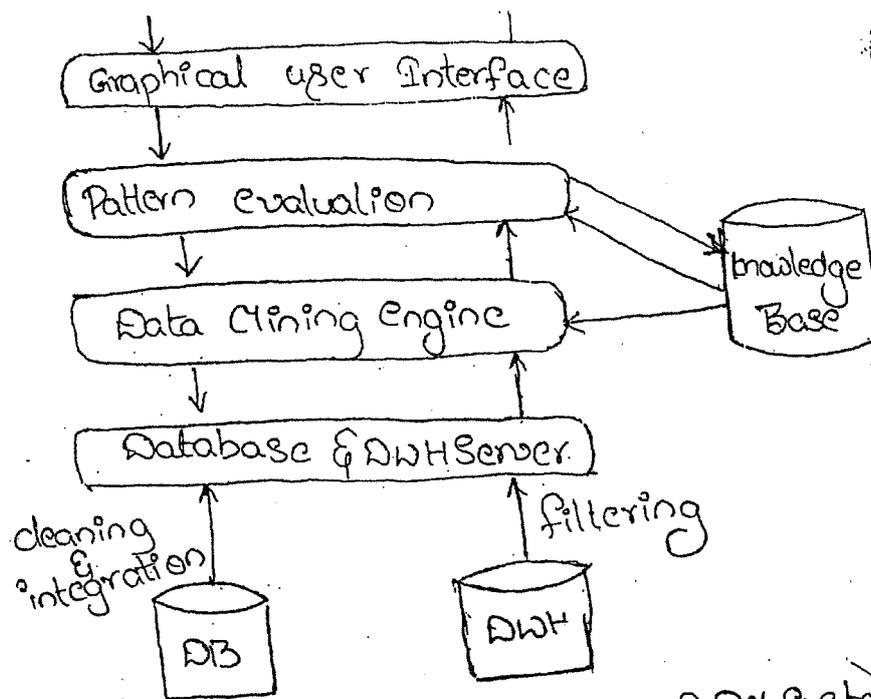
Here, DCI Systems extract the essential features from db or DWH. These Systems extract the data from centralized depository & the results are send to any file or db or DWH. These Systems are better than the no. of coupling systems.

3. Semitight Coupling:

Here, DCI Systems integrate with db or DWH to get the DCI primitives. i.e., store, retrieve, histogram analysis, join indexing etc.,

4. Tight Coupling:

Here, DCI Systems fully integrated with db or DWH to get all the features & to provide the integrated information environment. This is the desired architecture. Finally using this tight coupling we design the typical architecture of DCI.



Fig(a): Typical Architecture of DM System

The typical architecture of DM System contains the following components:

① DB, DWH (or) Any information Repository:

Here, we extract the data from any db (or) DWH or any information Repository then we apply the data cleaning, data transformation, data integration & finally data load. This data is loaded into db or DWH server.

② DB (or) DWH server:-

It contains the data according to user specification.

④ Knowledge Base:

Here, end user can get the knowledge from the different data sources. If the end user use the knowledge then the DSI process is simplified.

⑤ Data mining engine:

Here, we apply the DSI functions.

⑥ Pattern evolution:

Here, interesting patterns are evaluated.

⑦ GUI:

It provides the communication b/w end user & DSI system. i.e., through this end user present the queries to the DSI system, and also discovered patterns are visualized to the end user by using different visualization techniques.

Concept Description: characterization

& comparisons

The DSI is mainly classified into

- two
1. Descriptive mining
 2. Predictive mining

The descriptive mining analyzes the collected data set in the form of summarization. i.e., it explains the general characteristics of data.

The Predictive mining analyzes the data in order to construct one or more models using these models we can predict the behaviour of a new data set.

The descriptive mining is called as Concept description.

What is Concept description :-

The concept is nothing but the collection of data. For eg., frequently-buys, graduate-students etc., The concept description summarizes and explains the general characteristics of data. The concept description is also called as class description. It contains the two types.

1. class characterization

2. class comparison

The class characterization explains the characteristics of collected data.

The class comparison compares the two or more collections of data.

Comparison b/w OLAP & Concept Descriptions

a) Aggregate functions vs Complex data.

The DWH & OLAP systems the data is stored in the form of multi dimensional data model. i.e., Data cubes. The data cubes contains the dimensions & measures. The measures are analysed by using the aggregate functions, i.e., Sum(), Avg() etc., But the Concept description in DB handle the complex data types like non numeric, text, special, multimedia etc.,

b) Manual vs automation.

The DWH and OLAP systems we have to specify the dimensions & also we have to specify the OLAP operations. But the Concept description in db we select the dimensions & also we select the data at multiple levels.

Data Generalization & Summarization Based Characterization.

For eg., the data is store in large db. then, this data generalization provides the data abstraction from low

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Conceptual level to high Conceptual
The data generalization mainly classified
into two .

- (1) Data cube approach- we have to see the 2nd chapter .
- (2) Attribute Oriented Induction (AOI) Approach .

② Attribute Oriented Induction (AOI) Approach:

This is introduced in the year of 1989 . This is the best Approach for data generalization & summarization based characterization . It contains the two steps .

- 1. Select Task Relevant data .
- 2. Data Generalization .

1. Select Task Relevant data:

for eg... end user analyse the characteristics of graduate students from the Big university - DB & the attributes are name, gender, major, Birth-Place, Birth-date, Residence & Phone number .

Note: The data cube approach is based on DWH Orientation . But the AOI approach based on Relational Database .

on, this is specified in DDL

Syntax -

```
use Big-university - DB
mine characteristics as "science-students"
in relevance to name, gender, major,
birth-place, birth-date,
residence, phone #
from student
where status in "graduates"
```

This is transformed into relation query

Syntax :

```
use Big-university - DB
select name, gender, major, birth-place, birth-date,
residence, phone #.
from student
where status in {"MBA", "MA", "Mcom"}
```

Once this query is executed, we get the table. This table is called as Task relevant data or initial working relation. This is shown in below.

name	gender	major	birth-place	birth-date	residence	phone #
John	M	Science	Hyd, AP India	12/07/76	123 Bhill Hyd	234567
Lisa	F	Engineering	vancouver BC, Canada	10/08/70	Mainst vanco -over	99784

Data Generalization

Again it contains two steps.

- ① Attribute Removal
- ② Attribute Generalization

① Attribute Removal

It contains the two rules.

Rule 1: If the attribute has large no of distinct values and no hierarchy is defined then remove that attribute from the relation.

Rule 2: If the attribute has high concept hierarchy is defined then remove the low concept hierarchy & replaced with high concept hierarchy.

② Attribute Generalization

If the attribute has the large no of distinct values & concept hierarchy is defined then we select the generalization threshold value. After applying the ACI rules in Table 1.1 we get the following steps.

① name: If contains the large no of distinct values & no concept hierarchy is defined. ∴ Remove that attribute from the working relation.

② Gender: It contains 2 values and it remains the same.

③ major: It contains 3 values i.e., Science engineering, & business. It remains

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the same.

④ birth-Place :

It's concept hierarchy is defined city, state & country then this is replaced with birth country

⑤ birth-date :

This attribute is transformed into age & age is transformed into age-range.

⑥ Residence : It's concept hierarchy is defined street to city. This attribute is replaced with residence-city.

⑦ Phone # : It contains the large no of distinct values & no concept hierarchy is defined.

∴ remove it from the relation.

Here, similar tuples are merged & specified by using the count attribute

Then, finally we get the table 1.2

gender	major	birth-country	age-range	residence-city	count
M	Science	India	20...30	Hyd	20
F	Engineering	Canada	30...40	Vancouver	30

Table (1.2) After applying AOI Rules, the final Relational Table.

Efficient Implementation of Attribute-oriented Induction.

For efficient implementation of AOI we use the following algorithm.

Algorithm :- Attribute-oriented Induction.

Input :- (i) DB, a Relational db

(ii) Databases, a DQ Query.

(iii) A-list, list of Attributes

(iv) Gen(a_i), It is the Generalization operators for each a_i .

(v) A-thresh. value (a_i), It is the Generalization Threshold value for each a_i .

Output :- P ← Prime relation

methods :- w ← task-relevant data (DB-Query)

Here 'w' is the initial working relation

1) P ← prepare for generalization (w)

(a) If the attribute a_i contains the large no of distinct values & no concept hierarchy is defined then remove that attribute from the working relation

(b) If the attribute ' a_i ' high level concept hierarchy is defined then remove the low level concept hierarchy & replace with high level concept hierarchy.

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Q. P ← Generalization(θ);

- (a) If the attribute contains the large no of distinct values & concept hierarchy is defined, then we select the Threshold range.
- (b) If the generalization contains the similar tuples these are merged by using the count attribute.

Simple θ Algorithm.

* Presentation of Derived Generalization:

we apply the θ on Relational DB then we get the generalization Relation. This is presented at end user by using several visualization techniques i.e., cross table, Bar charts, Pie charts etc.,

For eg., Consider the sales table for all electronics is shown in below.

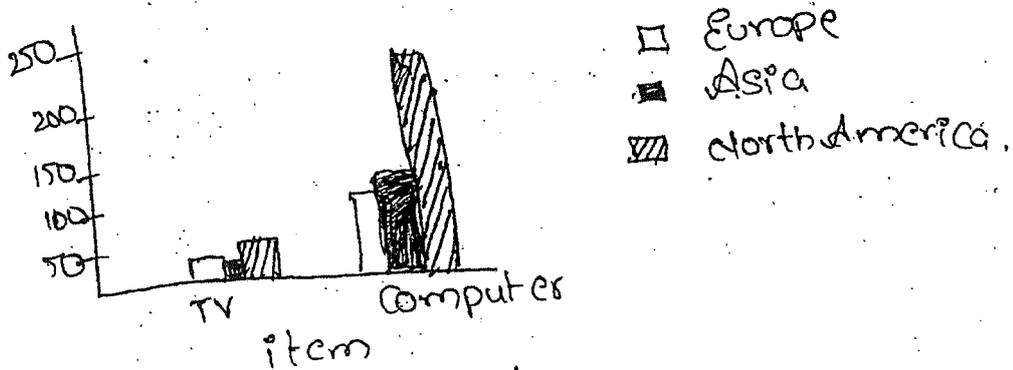
location	item	Sales (in dollars)	count (in thousands)
Europe	TV	15	200
Asia	TV	12	350
North America	TV	28	400
Europe	Computer	120	1000
Asia	"	130	1200
North America	"	250	1800

Result: Sales of all electronics

This table information is represented in cross table.

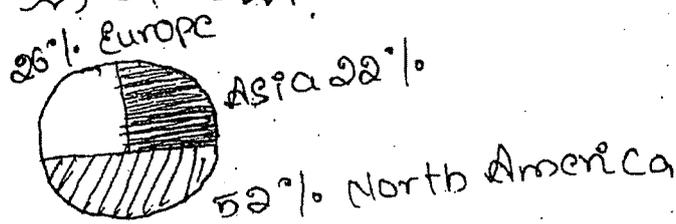
location/item	TV		Computer	
	Sales	Count	Sales	Count
Europe	15	200	120	1000
Asia	12	350	130	1200
North America	18	400	250	1800

fig(b) - cross table.

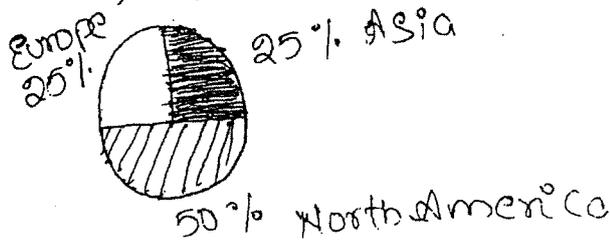


fig(c) Bar chart.

Pie chart for TV Sales:



Pie chart for Computer Sales:



* Analytical Characterization :-

using this attribute relevance we find the weakly relevant & irrelevant attributes. These are removed from concept description. using this we also find the most relevant attributes these are included in concept description.

why attribute relevance is acquired :-

The DWH & OLAP systems contains the drawbacks of enduser has to specify the dimensions & also he has to specify the high conceptual level. This is specified by using stmt "generalize dimension location to the country level".

The enduser does not know the attributes required to achieve the sql task then he specify all the attributes by using the stmt in relevance to *. But, this does not give the accurate data. To identify mostly relevant attributes & delete weakly relevant attributes we acquire the attribute relevance analysis.

Methods for Attribute Relevance Analysis :-

In this method we integrate the information gain analysis with

data dimension analysis

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finally, we find the information gain for each attribute & then the highest information gain attribute is added in concept description & remove the lowest information gain attribute from the concept description.

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for eg., Consider the training samples 's' & also each training sample class-label must be noted. To identify the class label we use one attribute for eg., Consider the attribute status using this attribute we find whether he is graduate student or under graduate student.

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lysis

let us consider the 'm' classes the sample 's' containing 's_i' samples in class 'c_i' for i=1, 2, ..., m then the sample belongs to class 'c_i' with the probability $\frac{s_i}{s}$ where s - total no of samples then the expected information needed to classify the given sample,

$$I(s_1, \dots, s_m) = - \sum_{i=1}^m \frac{s_i}{s} \log_2 \frac{s_i}{s} \quad \text{--- (1)}$$

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let us consider the attribute partition values $\{a_1, a_2, \dots, a_v\}$ then th

attribute position the given set 'S' into $\{s_1, \dots, s_m\}$ then let us consider the s_j falling C_{ij} then entropy of A.

$$E(A) = \sum_{j=1}^m \frac{(s_{1j} + \dots + s_{mj})}{S} \times I(s_{1j}, \dots, s_{mj}) \quad \text{--- (2)}$$

finally, information gain(A) = $I(s_1, \dots, s_m) - E(A)$

This method contains the following steps.

1. Data Collection :-

we collect the data for target class & contrasting class. The target class contains the data that is to be characterized. The contrasting class contains the ^{comparable} data that is to be characterized. ~~The~~

2. Apply AOI Rules :-

Here, we apply the AOI Rules.

3. Remove irrelevant & weakly Relevant Attributes :-

It contains three steps

Step 1 :- Compute the expected information gain

Step 2 :- Compute the entropy value.

Step 3 :- Compute the information gain. The highest information gain attributes is added in Concept description.

Example for class characterization :-

It contains the following steps

Step 1 :- collect the data for target class. i.e.,



to

graduate students. This is shown in below.

gender	major	birth-country	age-range	Count
M	Science	CANADA	20.....25	24
F	Science	CANADA	20.....25	30
M	Science	USA	20.....25	30
M	Engineer	CANADA	25.....30	20
F	Engineer	USA	25.....30	16

fig(a) Data for target class i.e., graduate students.

Step 2:- collect the data for contrasting class i.e., under graduate students.

gender	major	birth-country	age-range	Count
M	Science	CANADA	20.....25	20
F	Science	USA	20.....25	28
M	Engineer	CANADA	25.....30	20
F	Engineer	USA	25.....30	26
M	Business	CANADA	30.....35	20
F	Business	USA	30.....35	22

fig(b) Data for Contrasting class i.e., under graduate students.

We already apply the AOF Rules & Remove the name, phone # because they contains the large no of distinct values & no concept hierarchy is defined.

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The graduate student class is represented by S_1 & under graduate student class is represented by S_2 . The S_1 contains the 120 tuples & S_2 contains the 130 tuples. \therefore Expected information needed $I(S_1, S_2)$

$$I(S_1, S_2) = -\frac{120}{250} \log_2 \frac{120}{250} - \frac{130}{250} \log_2 \frac{130}{250}$$

$$= 0.99$$

Then consider the attribute major, for example, major = "Science".

$S_{11} = 84$, $S_{21} = 42$ then total = 126
expected information needed

$$I(S_{11}, S_{21}) = -\frac{84}{126} \log_2 \frac{84}{126} - \frac{42}{126} \log_2 \frac{42}{126}$$

$$= 0.78$$

major = "Engineer".

$S_{12} = 36$, $S_{22} = 46$ then total = 82

expected information needed

$$I(S_{12}, S_{22}) = -\frac{36}{82} \log_2 \frac{36}{82} - \frac{46}{82} \log_2 \frac{46}{82}$$

major = "Business"

$S_{13} = 0$, $S_{23} = 42$ then

expected information needed.

$$I(S_{13}, S_{23}) = 0$$

Then, entropy value for major

$$E(\text{major}) = \frac{126}{250} \times I(S_{11}, S_{21}) + \frac{82}{250} I(S_{12}, S_{22}) +$$

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$$\frac{42}{250} I(S_1, S_2)$$

to-
ced

$$= \frac{126}{250} \times 0.78 + \frac{82}{250} \times 0.98 + \frac{42}{250} \times 0$$

$$= 0.77$$

∴ information gain for major

$$G(\text{major}) = I(S_1, S_2) - E(\text{major})$$

$$= 0.99 - 0.77 = 0.22$$

ple...

Similarly, we can find the information gain for other attributes.

$$\text{e.g., } G(\text{gender}) = 0.003$$

$$G(\text{birth-country}) = 0.004$$

$$G(\text{age-range}) = 0.59$$

For e.g., Consider the attribute-relevance

Threshold is "0.1" then less than or equal to this value. Consider as the weakly relevant attributes.

∴ The attributes major & age-range is considered as the most relevant attributes these are included in Concept description. Similarly, the attributes gender & birth-country are the weakly relevant attributes these are removed from Concept description.

* Class Comparison : Discrimination by different class :-

The end user always have interest in mining the data from one or more contrasting

Classes. This is called as class comparison or class discrimination.

Rather than extracting the data from single class. i.e., class characterization. The target class & contrasting class must have the similarities. For eg., Person & item does not comparable. But the sales of last 3 years all comparable & also graduate students vs under graduate students are comparable.

* Methods for class comparison:

It contains the following steps.

Step 1: - collect the data for target class & contrasting class through DB queries.

Step 2: Apply the attribute relevance analysis.

Step 3: Apply the Generalization. i.e., we select the generalization threshold range for specific attribute. Then this is applied on target class & contrasting class to get the prime relations.

Step 4: - The derived class comparisons are presented to end user by using several visualization techniques.

Step 1: for eg., consider the big-university DB & the attributes are name, gender,

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major, birth-place, birth-date, residence & phone # to get the

use Big-university-DB

time Comparison as "graduate-vs-undergraduate-students"

in relevance to name, gender, major, birth-place, birthdate, residence, phone #

for "graduate-students"

where status in "graduate"

versus "undergraduate-students"

where status in "undergraduate"

analyse count %

from student.

Once the query is executed, we get the data for target class i.e., graduate students.

Name	gender	major	birth-place	birth-date	residence	phor
Mahi	M	Science	Hyd, AP INDIA	12-07-76	st-stres Hyd	8345
Lee	F	Engineering	Vancouver ABC, CANADA	12-07-70	123-st Vancouver	599-E

Fig(a) Data for Target class i.e., graduate student

Name	gender	major	birth-place	birth-date	residence	Phone#
Ravi	M	Science	N.Y. ABC, USA	12-07-80	Main St, N.Y	123456
Anu	F	busine -EE	Vancouver XYZ, CANADA	10-09-77	SR-st, Vancouver	234-578

Part(b) Data for contrasting class i.e., under grad. students

Step 2: Apply the attribute relevance analysis after applying the attribute relevance analysis the attribute name, gender, birth-place, residence & Phone # are removed from the relational table.

Step 3: Apply the generalisation. i.e., the attribute birth-place is transformed into age-range. The Prime relations are shown in below table.

major	age-range	count %
Science	20.....25	5.32%
Engineering	25.....30	4.12%

Fig(a) Prime Relation for graduate Students

major	age-range	Count %
Science	20.....25	3.32%
Business	30.....35	2.14%

Fig(b) Prime Relation for undergraduate Students

Step 4:

These Prime relations are presented to end user by using visualization techniques. In visualization technique we analyzed the

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measure is count %. for eg., 5.32% of students with the age 20.....25 & major in Science.

Presentation of class Comparisons:-

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2.

To Present the class Comparisons we use the method. This method is called as Statistics discrimination (or) d-weight value.

for eg., let s_A be the any tuple of the target class C_j then d-weight value of s_A in C_j is defined as the tuples containing target class by using total no of tuples.

$$d\text{-weight} = \text{count}(s_A \in C_j) / \sum_{i=1}^m \text{count}(s_A \in C_i)$$

Here, m contains the total no of tuples i.e., target class & contrasting class. for eg., consider the data for graduate students and under graduate students.

Status	major	age range	count
graduate	Science	20.....25	210
undergraduate	Engineering	25.....30	90

fig(a): Data for graduate & undergraduate students

The d-weight value for target class i.e., graduate students

$$d\text{-weight} = \frac{210}{300} = 0.7 = 70\%$$

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d-weight value for contrasting class.

i.e., undergraduate students

$$d\text{-weight} = 90/300 = 0.3 = 30\%$$

finally, if the student is major in science & with age 20...25 then, there is a probability is that he is the 70% of graduate students.

* Class Description:-

The class description means it contains the 2. i.e., class characterization & class comparison. for eg., Consider the data shown below.

Location	TV	Computers
Europe	90	150
North America	210	520

Fig(a): class table data for 2 locations

Then d-weight value i.e., TV

$$\text{Europe (d-weight)} = 90/300 = 0.3 = 30\%$$

$$\text{North America (d-weight)} = 210/300 = 0.7 = 70\%$$

d-weight value i.e., Computer

$$\text{Europe (d-weight)} = 150/670 = 0.22 = 22\%$$

$$\text{North America (d-weight)} = 520/670 = 0.78 = 78\%$$



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These d-weight values are shown in below.

Location / Item	TV		Computer	
	count	d-weight	count	d-weight
Europe	90	30%	150	22%
North America	210	70%	520	78%

Fig (b) d-weight value for fig(a)

from the fig(b) we can conclude that the north America sales are very good ✓

* Clining Statistical class Description from the large Database :-

upto now we analyze the characteristics of a relational DB by using the predefined aggregate functions like sum(), count() etc. But many of the applications end user want to analyze the characteristics of a large DB by using the measures mean, median, mode. These 3 measures are called as "Central tendency".

* Meas using the Central tendency :-

The most popular method to find the Centre value of a given set is mean. for eg...
Data items x_1, \dots, x_n then mean = $\frac{1}{n} \sum_{i=1}^n x_i$

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=78%

This mean is equal to avg in SQL.

$$\text{average} = \frac{\text{sum}}{\text{count}}$$

Sometimes the data contains the bytes - i.e., w_i for $i=1$ to n . Then, weighted arithmetic mean.

$$= \frac{\sum_{i=1}^n x_i w_i}{\sum_{i=1}^n w_i}$$

We can find the median value by using a formula:

$$\text{median} = L_1 + \left(\frac{N/2 - (\sum f)_1}{f_{\text{median}}} \right) \cdot c$$

where L_1 is the lowest class boundary
 c is the class interval.

$(\sum f)_1$ is the cumulative frequency $< f_{\text{median}}$

f_{median} is the median frequency

N is the no of samples.

Eg:- for eg., Consider the below table.

class interval 'c'	frequency f_{median}	Cumulative frequency $(\sum f)_1$
1-3	2	2
4-6	3	5
7-9	5	10

6. Mining Association Rules in Large Data Bases

Association Rule Mining:

The association rule Mining searches the interesting relationships between items in the given data set.

An typical example of association rule of Market Basket Analysis.

Market Basket Analysis:

This Market basket analysis analyse buying behaviour of customer by using association rule between the items.

For eg: Consider the three customers, then Market Basket Analysis is shown below.

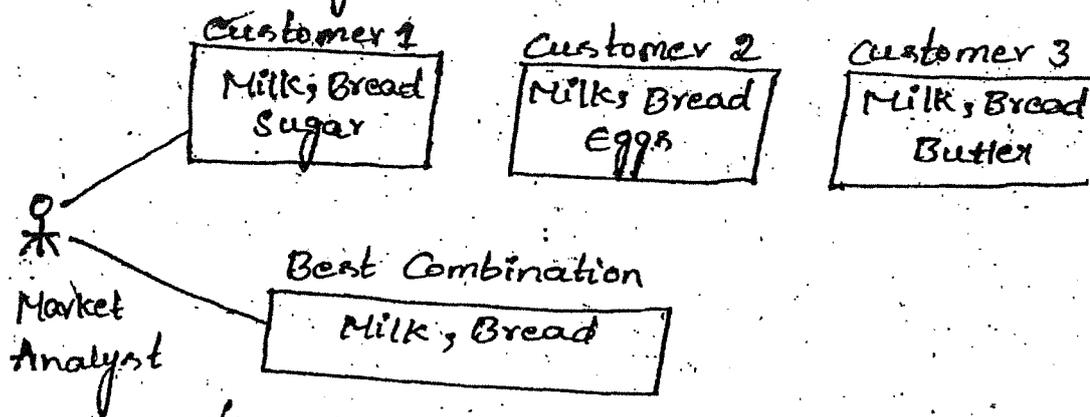


Fig: Market Basket Analysis.

Here Market Analyst identifies the customer purchases milk & also Bread. Therefore this milk and Bread placed together further more increase sales.

Therefore, as well as using the results of Market Basket Analysis the

1. The first part of the document discusses the importance of maintaining accurate records of all transactions.

2. It is essential to ensure that all entries are supported by proper documentation and receipts.

3. Regular audits should be conducted to verify the accuracy of the records and identify any discrepancies.

4. The final section outlines the procedures for handling any errors or corrections that may arise.

This market Basket Analysis uses the association rule.

For eg: Customer purchases a Computer and also purchases financial Mgmt Software. Then the association rule,

Computer \Rightarrow financial - Management - Software
[Support 2% ; Confidence = 60%]

Here Confidence 60% means 60% of customers purchase Computers & also purchase financial Mgmt Software & Support is 2%.

Basic Concepts:-

Consider Set of items, $I = \{i_1, i_2, \dots, i_m\}$ & task relevant data Set 'D'. we get this task relevant data set by using set of transactions. Each transaction is represented by $T_i, T \subseteq I$.

As well as each transaction is uniquely identified by using TID [Transaction ID].

For eg, item ACT, then $A \Rightarrow B, ACI, BCI$ & $A \cap B = \emptyset$

Then association rule $A \Rightarrow B$ with Support 's', where 's' is the percentage of transactions in 'D' contains A & B i.e.,

$$D(A \cup B)$$

The association rule $A \Rightarrow B$ has Confidence 'c', where 'c' is the percentage of transactions in 'D' contains A & also contains B i.e.,

ation

$$P(B/A)$$

\therefore Support $(A \Rightarrow B) = P(A \cup B)$ i.e., Transaction Contains 'A' & 'B'.

$$\text{Confidence } (A \Rightarrow B) = P(B/A) \text{ i.e.,}$$

Transaction Contains 'A' & also contains 'B'.

Here, Support & Confidence is represented in the form of percentage & ranges in between 0% to 100%.

Association Rule Mining : A Road Map :-

The market - Basket analysis i.e., one of the association rule. The association rule is classified based on the following criteria.

i. Based on the type of value used in rule :-

The association rule does not contain any item or attribute or dimension, then that association rule is called as boolean association rule.

Ex: Computer \Rightarrow financial - management - Software

If the association rule is described by quantitative values, then that association rule is called as quantitative association rule.

Ex: $\text{age}(x, "20 \text{ --- } 29") \wedge \text{income}(x, "30k \text{ --- } 39k")$
 $\Rightarrow \text{buys}(x, "VCR")$

ii. Based on the type of dimensions used in the :

For Ex: The association is defined for one

dimension then that association rule is called as single dimension association rule.

Ex: $\text{buys}(x, \text{"computer"}) \Rightarrow \text{buys}(x, \text{"financial-Mgmt-S/w"})$

iii) Based on the different types of Abstractions
in the Rule:

Here the association rule contains the different types of data abstractions. i.e. using this association rule we abstract data at different levels.

Here the association rule contains the different types of data levels.

Ex: $\text{age}(x, \text{"30 --- 39"}) \Rightarrow \text{buys}(\text{"Computer"})$

$\text{age}(x, \text{"30 --- 39"}) \Rightarrow \text{buys}(\text{"laptops"})$

** Mining Single dimensional Boolean Association Rule From Transactional Dbs:

Here we use the 'Apriori' algorithm.

Apriori Algorithm: Finding frequency Itemsets
by using candidate sets.

- 1) The Apriori Algorithm is used to find frequent Item sets.
- 2) These are used in Boolean association rule.
- 3) Apriori means recursive approach, i.e., level-wise-search.

4) First of all we find frequent itemset,
This is denoted by L_1 . Then using L_1
we find L_2 i.e., frequent 2-itemset.

5) This is used to find L_3 and so on,
until we can not get the no frequent set i.e.,
the Apriori algorithm uses prior knowledge
of frequent item set and for each L_k we have
to scan entire database i.e., D .

To simplify the Apriori algorithm, it
uses apriori property. Using this Apriori
Property, we reduce search length.

For ex, item set is represented by 'I' & it
does not satisfying the minimum support
threshold i.e., sup-min , then I is not a
frequent itemset, then

$$P(I) < \text{sup-min}$$

If we add another item 'A' then it results
IVA,

It is also not a frequent itemset

$$P(P(I)) < \text{sup-min}$$

This apriori property contains the term steps

1) The Join Step:

To calculate L_k , a candidate k itemset
are joined, L_{k-1} to itself. The candidate set
is represented by C_k . Consider L_1 & L_2 item

in l_{k-1} . In general $x_1 \leq x_2$ means x_1 value of l_{k-1} , then we join $l_{k-1} \times l_{k-1}$, it requires the first $k-2$ items are equal. Therefore, to join l_1 and l_2 of l_{k-1} and if

$$[l_1[1]=l_2[1]] \wedge [l_1[2]=l_2[2]] \wedge \dots \wedge [l_1[k-2]=l_2[k-2]]$$

$$\wedge [l_1[k-1] < l_2[k-1]]$$

The last condition,

$l_1[k-1] < l_2[k-1]$ is used to avoid the duplicate values.

2) The Prune Step:-

The prune step, all non-empty subset of frequent items is also frequent.

Ex: Consider transactional database of 'all electronics'. This is represented by 'D' & it contains of transactions.

\therefore we use the Apriori algorithm to find frequent item & have minimum count is '2' C_1

TID	list of item-IDs	Scan D for count of candidates	item set	Support Count
T100	I1, I2, I5	→	{I1}	6
T200	I2, I4		{I2}	7
T300	I2, I3		{I3}	6
T400	I1, I2, I4		{I4}	2
T500	I1, I3		{I5}	2
T600	I2, I3			
T700	I1, I3			
T800	I1, I2, I3, I5			

High transaction per an customer

Compare Candidate Sup-count with min sup-count

d_1

Item set	Sup-count
{I1}	6
{I2}	7
{I3}	6
{I4}	2
{I5}	2

$C_2 = d_1 \times d_1$

Generate C_2 Candidate using d_1

Itemset	Sup-count
{I1, I2}	4
{I1, I3}	4
{I1, I4}	1
{I1, I5}	2
{I2, I3}	4
{I2, I4}	2
{I2, I5}	2
{I3, I4}	0
{I3, I5}	1
{I4, I5}	0

Compare Candidate Sup-count with min sup-count

d_2

Itemset	Sup-co
{I1, I2}	4
{I1, I3}	4
{I1, I5}	2
{I2, I3}	4
{I2, I4}	2
{I2, I5}	2

$C_3 = d_2 \times d_2$

Generate C_3 Candidate using d_2

Itemset	Supcount
{I1, I2, I3}	2
{I1, I2, I5}	2

Compare Candidate C-c sup-count with min sup-count

d_3

itemset	S_c
{I1, I2, I3}	
{I1, I2, I5}	

Using Apriori algorithm, we find candidate & most frequently item set with min Support Count

Steps for finding candidate item set of frequent items

1) This algorithm scans all the transactions of database to find total count of each item set. This item set is called as candidate 1-itemset, C_1 .

2) Suppose consider minimum support threshold is 2 ($2/9 = 0.22 = 22\%$)

3) Candidate Support Count is compared with minimum support count. Then we get frequent itemset d_1 .

4) To find C_2 , the d_1 joined itself i.e.

$$\text{Join } C_2 = d_1 \bowtie d_1$$

Then we find d_2 by comparing candidate support count with minimum support-count.

5) Join $C_3 = d_2 \bowtie d_2$

$$\begin{aligned} & \{ \{I_1, I_2\}, \{I_1, I_3\}, \{I_1, I_5\}, \{I_2, I_3\}, \{I_2, I_4\}, \\ & \quad \{I_2, I_5\} \} \bowtie \{ \{I_1, I_2\}, \{I_1, I_3\}, \{I_1, I_5\}, \\ & \quad \{I_2, I_3\}, \{I_2, I_4\}, \{I_2, I_5\} \} \\ = & \{ \{I_1, I_2, I_3\}, \{I_1, I_2, I_5\}, \{I_1, I_3, I_5\}, \\ & \quad \{I_2, I_3, I_4\}, \{I_2, I_3, I_5\}, \{I_2, I_4, I_5\} \}. \end{aligned}$$

Prune Step:-

All non-empty frequent subsets are also frequent.

1. 2-Item sets for $\{I_1, I_2, I_3\}$ are $\{I_1, I_2\}$, $\{I_1, I_3\}$ and $\{I_2, I_3\}$. Then $\{I_1, I_2, I_3\}$ Add to C_3 because all 2-item sets present in d_2 .
2. 2-Item sets for $\{I_1, I_2, I_5\}$ are $\{I_1, I_2\}$, $\{I_2, I_5\}$ and $\{I_1, I_5\}$ - there all 2-item set Present in d_2 Then $\{I_1, I_2, I_5\}$ add to C_3 .
3. 2-Item sets for $\{I_1, I_3, I_5\}$ are $\{I_1, I_3\}$, $\{I_3, I_5\}$ and $\{I_1, I_5\}$ Then $\{I_3, I_5\}$ is \cap in d_2 then $\{I_1, I_3, I_5\}$ is Removed from C_3 .
4. 2-Item sets for $\{I_2, I_3, I_4\}$ are $\{I_2, I_3\}$, $\{I_3, I_4\}$ and $\{I_2, I_4\}$ then $\{I_3, I_4\}$ is not in d_2 . Therefore, it is not a frequent set. So $\{I_2, I_3, I_4\}$ is removed from C_3 .
5. 2-Item sets for $\{I_2, I_3, I_5\}$ are $\{I_2, I_3\}$, $\{I_3, I_5\}$ and $\{I_2, I_5\}$ Then $\{I_2, I_3, I_5\}$ is removed from C_3 .
6. 2-Item sets for $\{I_2, I_4, I_5\}$ are $\{I_2, I_4\}$, $\{I_4, I_5\}$ and $\{I_2, I_5\}$ then $\{I_4, I_5\}$ is \cap in d_2 . Therefore it is not frequent set. Then $\{I_2, I_4, I_5\}$ is removed from C_3 .

7. Then $C_3 = \{ \{I_1, I_2, I_3\}, \{I_1, I_2, I_5\} \}$
8. Again we scan all the transactions to find the count value for 3-itemsets. Then we get d_3 .

Item set	Sup-count
$\{I_1, I_2, I_3\}$	2
$\{I_1, I_2, I_5\}$	2

Apriori Algorithm:

Find frequent set using recursive level-wise approach based on candidate item-set.

Input: Database D , Op-transaction; minimum Support threshold, min-sup

Output: d , frequent itemset.

Method: $d_1 = \text{first-frequent-1-itemset}$;

for ($k=2$; $d_{k-1} \neq \emptyset$; $k++$)

{

$C_k = \text{apriori-gen}(d_{k-1}, \text{min-sup})$;

for each transaction $L \in D$

{

$C_k = \text{Subset Count}$

for each $C \in C_k$

$C \cdot \text{count}++$;

}

$d_k = \{ C \in C_k \mid C \cdot \text{count} \geq \text{min-sup} \}$

```

}
return L;
}
Procedure apriori-gen ( $L_{k-1}$ : frequent  $k-1$  item set
min-sup: min. support threshold)
{
Consider  $l_1, l_2, l_1 \in L_{k-1}, l_2 \in L_{k-1}$ 
if  $l_1[1] = l_2[1] \wedge \dots \wedge (l_1[k-2] = l_2[k-2])$ 
 $(l_1[k-1] < l_2[k-1])$  then  $C = l_1 \bowtie l_2$ ;
if has-infrequent-subset ( $C$  vs  $L_{k-1}$ ) then
delete  $C$ ;
else
add  $C$  to  $C_k$ ;
return  $C_k$ ;
}

```

```

Procedure has-infrequent-subset ( $C$ : candidate
item set,  $L_{k-1}$ : frequent  $k$ -item set)
{

```

```

if  $C \notin L_{k-1}$  then
return TRUE
else
return FALSE
}

```

Alg:-

Apriori Algorithm for finding frequent item
i.e., used in boolean association rule.

Generating Association rules from frequent item sets

Using Apriori alg, we find frequent item sets then we derive strong association rules

The strong association rule means it must include Support Count & Confidence Count. Then we use the following equation for the Confidence to derive strong association rules where Conditional Probability is represented in the form of the Support Count.

$$\text{Confidence } (A \Rightarrow B) = P(B/A) = \frac{\text{Support-Count } (A \cup B)}{\text{Support-Count } (A)}$$

where Support-Count $(A \cup B)$ represents transactions containing A & B.

Support-Count (A) represents transaction containing 'A'.

Using this we derive 2 associations rules for each frequent item sets L , must contain all non-empty subsets of 'L'.

For each subset 'S' of 'L', the o/p rule

$$"S \Rightarrow (L-S)" \text{ if } \frac{\text{Support-Count}(L)}{\text{Support-Count}(S)} \geq \text{min-Confidence}$$

For ex, frequent itemsets,

$L = \{I_1, I_2, I_5\}$ then subsets are

$\{I_1, I_2\}$, $\{I_1, I_5\}$, $\{I_2, I_5\}$, $\{I_1\}$, $\{I_2\}$ & $\{I_5\}$

Then the association rules are

- $I_1 \cap I_2 \Rightarrow I_5$, Confidence = $2/4 = 50\%$ — ①
- $I_1 \cap I_5 \Rightarrow I_2$, Confidence = $2/2 = 100\%$ — ②
- $I_2 \cap I_5 \Rightarrow I_1$, Confidence = $2/2 = 100\%$ — ③
- $I_1 \Rightarrow I_2 \cap I_5$, Confidence = $2/6 = 33\%$ — ④
- $I_2 \Rightarrow I_1 \cap I_5$, Confidence = $2/7 = 28\%$ — ⑤
- $I_3 \Rightarrow I_1 \cap I_2$, Confidence = $2/2 = 100\%$ — ⑥

Consider min confidence is 70% , then the o/p is ②, ③, ⑥ only [we get] strong association

Improve the Efficiency of Apriori:-

we use following techniques to improve the efficiency of Apriori alg.

1. Hash based Technique:-

using this , we reduce the size of candidate - k itemset . we create both hash table for 'H' by using the hash function.

$$H(x, y) = (\text{Order of } x) \times 10 + (\text{Order of } y) \pmod 7$$

we consider transaction db G-1 then hash table ht₂ for Candidate - 2 itemset.

Hash address	0	1	2	3	4	5	6
Hash Count	2	2	4	2	2	4	4
Hash Contents	{I ₄ , I ₄ } {I ₂ , I ₅ }	{I ₁ , I ₅ } {I ₁ , I ₅ }	{I ₂ , I ₃ } {I ₂ , I ₃ } {I ₂ , I ₃ } {I ₂ , I ₃ }	{I ₂ , I ₄ } {I ₂ , I ₄ }	{I ₂ , I ₅ } {I ₂ , I ₅ }	{I ₁ , I ₂ } {I ₁ , I ₂ } {I ₁ , I ₂ }	{I ₁ , I ₂ } {I ₁ , I ₂ }

$$\{I_1, I_2, I_5\}$$

$$h(I_1, I_2) = 10 + 2 \cdot 7 = 5$$

$$h(I_1, I_5) = 10 + 5 \cdot 7 = 1$$

$$h(I_2, I_5) = 20 + 5 \cdot 7 = 4$$

$$h(I_2, I_4) = 20 + 4 \cdot 7 = 3$$

$$h(I_2, I_3) = 20 + 3 \cdot 7 = 2$$

$$h(I_1, I_4) = 10 + 4 \cdot 7 = 0$$

$$\{I_1, I_2, I_4\} \Rightarrow h(I_1, I_4) = 10 + 4 \cdot 7 = 0$$

$$h(I_1, I_3) = 10 + 3 \cdot 7 = 6$$

$$\{I_1, I_2, I_3, I_5\}$$

$$h(I_3, I_5) = 30 + 5 \cdot 7 = 0$$

Consider min-support count is '3'. Then the addresses 0, 1, 3, 4 does not be considered for c2.

2. Transaction Reduction:-

If the transaction does not contain frequent k-itemset, then it also does not contain the frequent k+1 itemset then those transactions must be removed.

3. Partitioning:-

In this, it requires only 2 database scans & also it consists 2 phases.

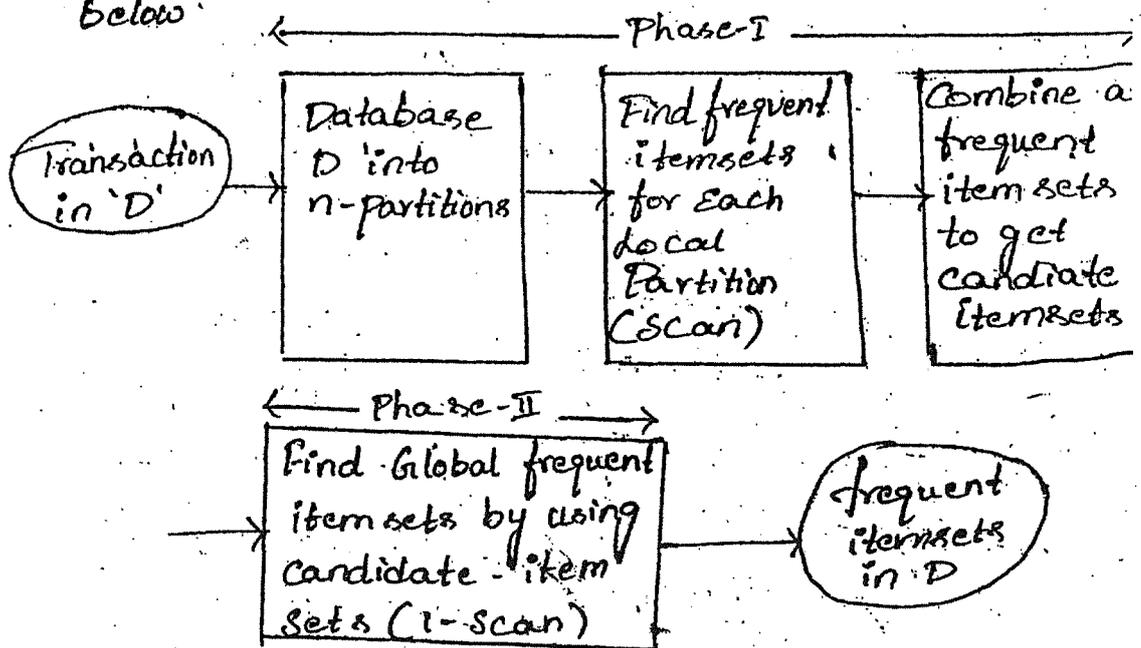
Phase-1: It contains following steps.

1. Divide db 'D' into 'n' partitions.
2. Find frequent itemset for each local partition

3. Combine these
Candidate item sets

Phase-II

1. Find global frequent item set - This is shown below



Partition data Method.

4. Sampling:

Consider the random sample 's' in database 'D'. Then we find frequent itemsets in 's' rather than 'D'.

The size of 's' is such a way that it can be fit in primary memory i.e., in single scan all the transactions of 's' are avoided.

Here we are finding frequent itemset in 's'. Therefore, we may miss some frequent itemsets. Therefore we consider lower support threshold rather than min. support threshold.

Generate Frequent itemsets without using Candidate itemsets (or) FP growth Method

In Apriori algorithm, we find Candidate itemsets. Using this Candidate itemsets, we find frequent itemsets. But the Apriori Algorithm contains the following drawbacks.

- i) It generates large No. of Candidate itemsets. For ex, 10^4 frequent-itemset. Then apriori alg generates approximately 10^8 Candidate-2 itemsets.
- ii) It requires the large No. of database Scans to identify the value of candidate itemsets.

To avoid these drawbacks, we go for the frequent pattern-growth Method.

This is simply called as FP-growth method. It follows the Divide-and-Conquer technique.

It contains three steps.

1. Compress the db representing itemsets into frequent pattern-tree or FP-tree. following the itemset information.
2. Transform Compressed db into conditional DB.
3. Generate frequent Patterns.

Consider the transaction db fig:

TID	list of item - l_k
T100	I_1, I_2, I_5
T200	I_2, I_4
T300	I_2, I_3
T400	I_1, I_2, I_4
T500	I_1, I_3
T600	I_2, I_3
T700	I_1, I_3
T800	I_1, I_2, I_3, I_5
T900	I_1, I_2, I_3

After the first scan of db, we get l_1 it contains the item set & support-count & we represent this in descending order of support count & it is represented by 'L'

$$L = [I_2:7, I_1:6, I_3:6, I_4:2, I_5:2]$$

Consider support-count is '2', then FP-tree construction is

after the second scan of db, the first transaction is

$$"T_{100} : I_1, I_2, I_5"$$

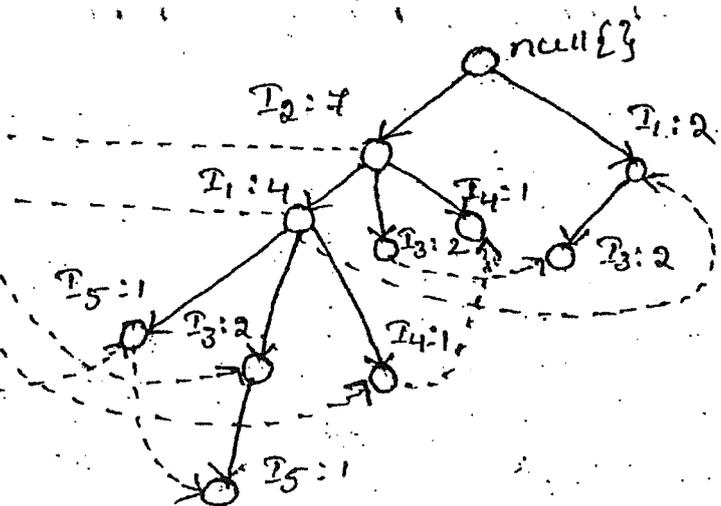
& these items are arranged based on the order of 'L'. Then we get (I_2, I_1, I_5) & these items are occurred in first time, the $(I_2:1)$ is treated as the root & $(I_1:1)$ is attached to I_2 & $(I_5:1)$ is attached to I_1 & then second transaction.

$T_{200} : I_2, I_4$

then, I_2 is already root, so increment the count value of ' I_2 ' & ($I_4:1$) is attached to I_2 . This is shown below.

✓ FP-Tree Construction Start with the root, this root is labelled as null { }

Item I_0	Support Count	Node LINK
I_2	7	
I_1	6	
I_3	6	
I_4	2	
I_5	2	



✓ FP-tree based on Transaction Database.

Construction of Conditional DB by Mining FP-Tree:

Start with the item ' I_5 ' it contains the 2 branches. These branches are identified by verifying Node link. Then the paths are

$I_2 I_1 I_5 : 1$ $I_2 I_1 I_3 I_5 : 1$

Then I_5 is considered as suffix. Then we

1. Conditional patterns how i.e., as link paths of I_5 are ($I_2 I_1 : 1$) and ($I_2 I_1 I_3 : 1$)

Then Conditional FP-tree Contains only one path, then we merge above 2 paths & we get

we cannot include I_3 because its support count is '1' & it is less than the support count. Then this single path generates the frequent itemsets by concatenating suffix.

∴ Frequent patterns are

$I_2 I_5 : 2$ $I_1 I_5 : 2$

This is shown below.

Item	conditional pattern Base	Conditional FP-tree	Frequent Pattern Generation
I_5	$\{ \langle I_2 I_1 : 1 \rangle, \langle I_2 I_1 I_3 : 1 \rangle \}$	$\langle I_2 : 2, I_1 : 2 \rangle$	$I_2 I_5 : 2$ $I_1 I_5 : 2$ $I_2 I_1 I_5 : 2$ $I_2 I_4 : 2$
I_4	$\{ \langle I_2 I_1 : 1 \rangle, \langle I_2 : 1 \rangle \}$	$\langle I_2 : 2 \rangle$	$I_2 I_4 : 2$
I_3	$\{ \langle I_2, I_1 : 2 \rangle, \langle I_2 : 2 \rangle, \langle I_1 : 2 \rangle \}$	$\langle I_2 : 4, I_1 : 2 \rangle, \langle I_1 : 2 \rangle$	$I_2 I_3 : 4, I_1 I_3 : 2, I_2 I_1 I_3 : 2$
I_1	$\{ \langle I_2 : 4 \rangle \}$	$\langle I_2 : 4 \rangle$	$I_2 I_1 : 4$

Conditional db by Mining FP-tree

Finally, using this FP-growth method large pattern is divided into smaller pattern by using suffix & then suffix is concatenated to get the frequency patterns.

This alg. reduces the search cost & also this is more faster than the apriori algorithm.

ICEBERG QUERIES:

The Apriori alg is used to improve the efficiency of ICE-BERG Query. This ice-berg query is mainly used in DM preferably Market

The ICE-BERG query computes aggregation function over an attribute, to find out which aggregate values over the specified threshold.

Consider the relation 'R' with attributes $a-1, a-2, \dots, a-n$, & b and aggregate function $agg-f$. Then the ice-berg query looks like the following.

Select $R.a-1, R.a-2, \dots, R.a-n, agg-f(R.b)$
from relation R.

group by $R.a-1, R.a-2, \dots, R.a-n$
having $agg-f(b) \geq \text{threshold}$.

The above ICE-BERG query accepts the large no of i/p tuples, but it gives very small No. of o/p tuples, because the o/p tuples only displays that satisfy the threshold values.

MINING MULTILEVEL ASSOCIATION RULES FROM TRANSACTION DATABASES:-

Mining multilevel association rules means rules are included at different levels of abstraction.

Multilevel Association Rules:-

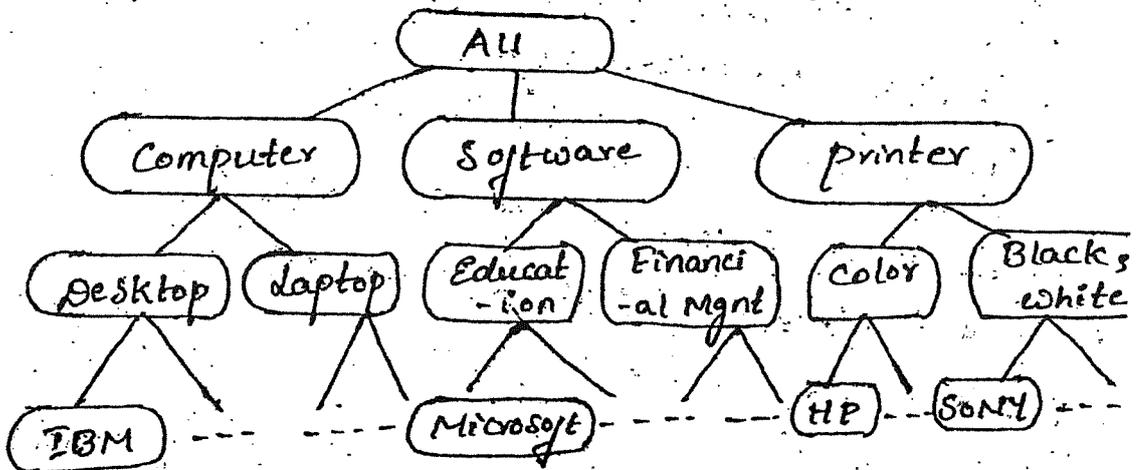
It is difficult to define strong association rules at low level or primitive level.

Consider the task-relevant data set as specified below.

TID	Items Purchased
T1	IBM, Desktop Computer, Sony, B/W printer,
T2	Microsoft Education S/w, HP Color Printer
T3	Microsoft financial Mgmt. S/w,
T4	IBM laptop Computer
⋮	⋮

Task Relevant Data set

The above task relevant data set is represented in Concept hierarchy as.



The Concept hierarchy defines set of mapping from low level to high level.

Here it contains 4 levels. i.e., levels 0, 1, 2, 3. The top level is level '0' i.e., represented by using keyword 'all'.

The level '1' contains Computer, S/w & Printer.

The level 2 contains Desktop Computer, Laptop Computer & so on.

The level 3 contains IBM Desktop Computer & so on.

The lowest level is level 0 & in this level it is difficult to identify interesting patterns i.e.,

1. Consider "IBM Desktop Computer" & "Sony B&W Printer" occur in very few.

∴ Define Generalization like "IBM Desktop Computers" to "Computer" & "Sony B&W printer" to "Printer".

∴ The combination {Computer, Printer} as many people frequently purchased.

∴ using this multilevel concept hierarchies, we easily define the interesting patterns.

Approaches to Mining Association Rules:

Generally we follow the top down approach, i.e., with level 1 & find frequent item

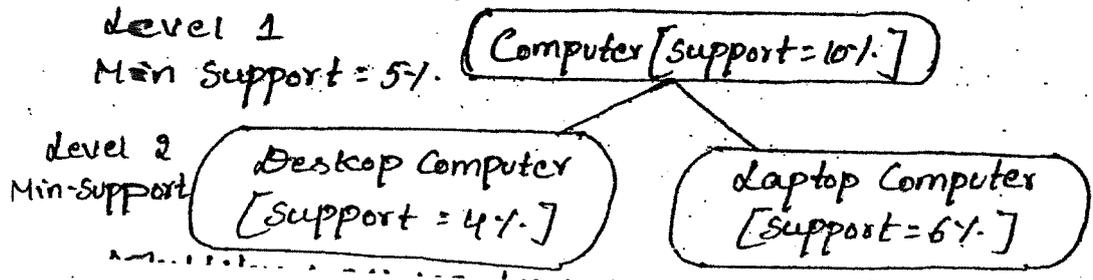
There are no frequent item sets. Then we go for level 2 & so on.

we use the following approaches for mining association rules.

Uniform Support:

Here we define same minimum support threshold for all the levels.

For α , Min. Support Threshold is 5%, then



Therefore, we can find frequent patterns & desktop Computer or not.

This approach contains the advantages of I. Searching process is simplified i.e., here we verify only those items satisfy Min-Support II. Here Enduser has to specify only one Min-Support Threshold Value.

This technique contains the drawbacks of

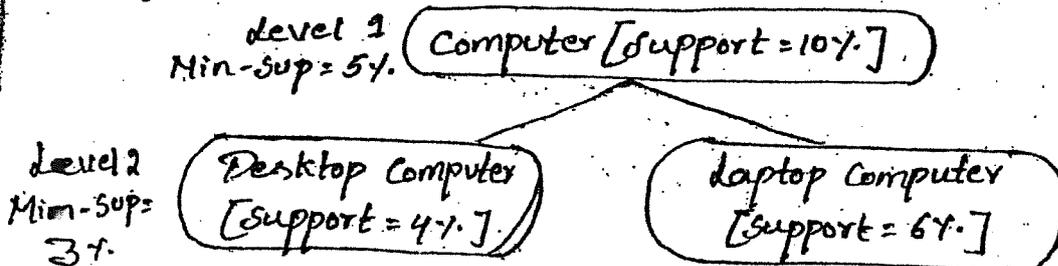
i) we already see the low level item is not frequent as & upper level item. Therefore, we cannot define Min-Support Threshold.

ii) if we specify very large support threshold then the low level important patterns may be missed.

iii) If we define very small support threshold then it contains the large No. of unimportant patterns.

ii) Reduced Support:

Here each level has its own minimum support threshold. Generally higher levels contains the big values than the low levels. Consider Min-Support threshold for level 1 & 2 are 5% & 3%. This is shown in below.



Multilevel Mining with reduced Support.

These computers, mainly computers, are frequent patterns.

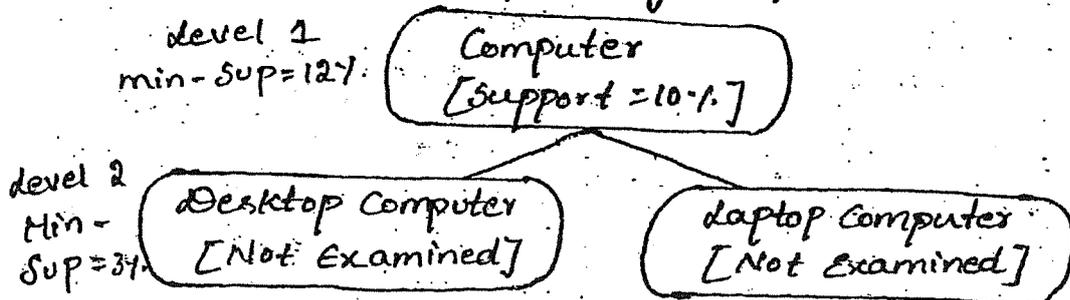
iii) level-by-level Independent :-

Here each node is examined independently i.e., regardless of its parent is frequent or not.

iv) level Cross Filtering by Single item :-

Here 'item with i th level is examined if & only iff its parent $(i-1)$ th level is frequent.

Consider the following Diagram.



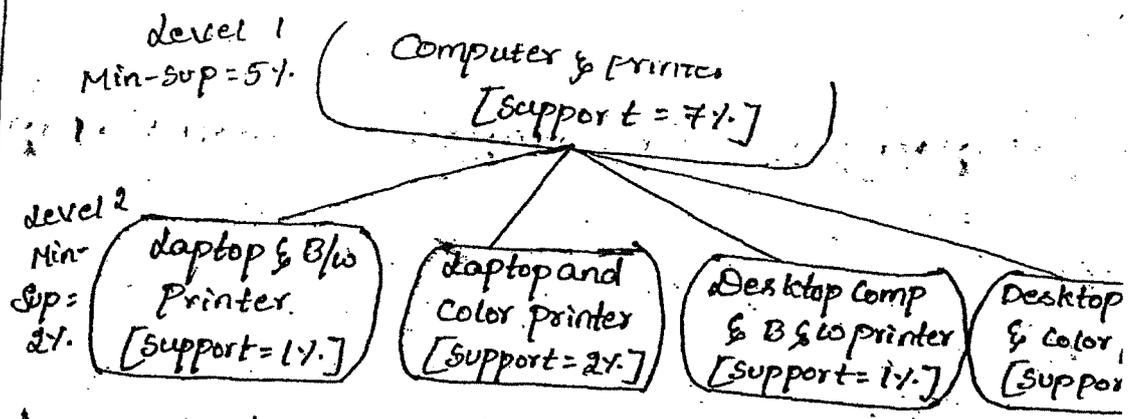
Multilevel Mining with reduced support using level cross filtering by single item.

Here Desktop Computer & Laptop Computer is not examined because its parent 'Computer' is not a frequent.

v) level Cross Filtering by k-item Set :-

Here k-item Set is examined [k-itemset with i th level] if & only iff Parent k-itemset of $(i-1)$ th level is frequent.

Consider the below diagram.



Multilevel Mining with Reduced Support using level-cross filtering by k-item set, here $k=2$

- Here { Laptop Computer, B/w Printer }
- { Laptop Computer, Color " }
- { " " , B/w " }
- { " " , Color " }

because 'Computer' & 'Printer' are frequent.

Checking redundant Multilevel Association Rule

using concept hierarchy we define association rules at different levels of abstraction. This is called as Multilevel association rule. when we are deriving multilevel association rules we may contain redundant rules. when we are deriving multilevel association rules we may contain redundant rules by "ancestor" property exist different items.

For Ex Consider the concept hierarchy fig 6.4 contains the ancestor property of 'Desktop Compu' is the ancestor of 'IBM'.

02

For Ex, Consider the following association rules.

Desktop Computer \Rightarrow B&W printer — ①

IBM " " \Rightarrow B&W printer — ②

Here rule ② doesn't give any ad

Let 'R₁' is the ancestor of 'R₂' then 'R₁' replaces 'R₂'.

Mining MultiDimensional Association rules from relational DB (or DWH Cor) Pattern Evolution Method:

— Here we mine association rules more than one dimension or predicates.

Multi Dimensional Association Rules:

For Ex, Consider the 'ALL ELECTRONICS' DB. The Boolean Association rule is i.e., it doesn't contain any dimension.

IBM Desktop Computer \Rightarrow Sony B&W Printer.

Using this we define single dimension association rule as

$buys(x, "IBM Desktop Computer") \Rightarrow$

$buys(x, "Sony B&W Printer").$

— Here we are using one predicate "buys" we can also define multidimensional Association rule as

$age(x, "20 --- 29") \wedge occupation(x, "Student")$

$\Rightarrow buys(x, "laptop").$

Here, it contains occupation & buys.

Here No. Predication is Repeated. Then it is called as Inter Dimensional Association Rule.

If a predicate is repeated, then called as hybrid Dimensional Association. This is shown below.

$age(x, "20-29") \wedge buys(x, "daptop") \Rightarrow buys(x, "B&W")$

Generally, attributes in database quantitative.

Here 'Categorical' means it contains finite no. values & this values does contain any ordering.

Quantitative means it contains any numerical values & this values are ordered.

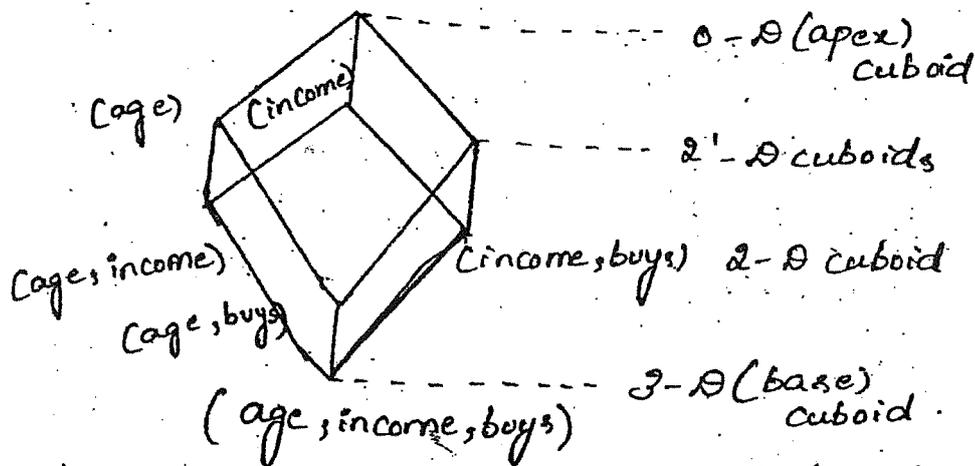
Mining Multidimensional Association using Stati.
Discretization of Quantitative:

Discretization is performed before the D.
Consider the concept hierarchy for in is replaced with numerical range of num values such as "20k --- 40k", "41k --- 80k".

Then the discretization is static & pre-defined.
This range of values consider as categories.
We find the values to fall in each category.
Then we find frequent item set using this freq item set, we derive association rules.

task relevant data cube. Data cube is a multidimensional structure that contains the multidimensions & also using the data cube, we easily derive multidimensional association rules.

The data cube consists of lattice of cuboids. These are multidimensional data structures & these contain task relevant data & also contain aggregated & grouping information. This is shown below.



Lattice of cuboids make 3-D Data cube. Here the base cuboid aggregates task relevant data by age, income, buys.

2-D cuboid i.e., (age, income), it aggregates task relevant data age & income.

1-D cuboid specifies task relevant data of 1 dimension.

The apex cuboid specifies total No. of transactions in task relevant data.

Mining Quantitative Association Rules

Quantitative association rules are multidimensional association rules but they contain numerical attributes. These are derived from discretization.

In Quantitative association rules, the left side it contains 2 Quantitative attributes & in right side it contains Categorical attributes. These is shown below:

$$A_{quan1} \wedge A_{quan2} \Rightarrow A_{cat}$$

Here, A_{quan1} , A_{quan2} are quantitative Numerical range values.

' A_{cat} ' is the categorical attribute task relevant data.

For Ex, 2-D Quantitative association rule is:
 $age(x, "30-39") \wedge income(x, "29k-30k") \Rightarrow buys(x, "TV")$

These quantitative association rules derived from ARCS (Association Rules Clustering System).

This ARCS approach contains the following steps.

1) Binning: The Quantitative attributes generally contain wide range of values. To smooth the values, we use following binning methods.

a) Equiwidth: here it contains interval size. Each bin is equal.

1) Κυματώδη: there is continuous change in values in each bin.

2) Find Frequent Set:

To find frequent set, we construct 2-dimensional array containing count values for each category. These values are scanned to get frequent set. These frequent set also satisfy min. support & min. confidence.

3) Clustering association rules:

For ex, consider the association rules below.

$$\text{age}(x, 34) \wedge \text{income}(x, "31k \dots 40k") \Rightarrow \text{buys}(x, "TV") \text{--- ①}$$

$$\text{age}(x, 35) \wedge \text{income}(x, "31k \dots 40k") \Rightarrow \text{buys}(x, "TV") \text{--- ②}$$

$$\text{age}(x, 34) \wedge \text{income}(x, "41k \dots 50k") \Rightarrow \text{buys}(x, "TV") \text{--- ③}$$

$$\text{age}(x, 35) \wedge \text{income}(x, "41k \dots 50k") \Rightarrow \text{buys}(x, "TV") \text{--- ④}$$

These rules are represented in 2D Grid as below

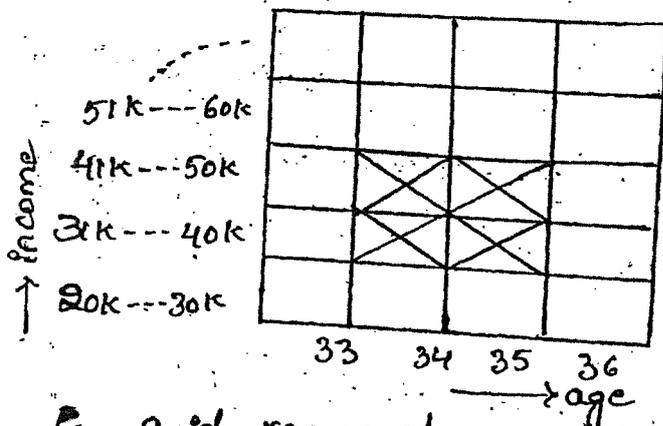


Fig: grid represents customer who buys "TV"

The above 4 association rules are simply represented in:

$$\text{age}(x, "34 \dots 35") \wedge \text{income}(x, "30k \dots 50k") \Rightarrow \text{buys}(x, "TV").$$

Mining Distance based

For ex, the data for the 'price' partitioned by using equiwidth & equidepth & this is compared with distance-based partition & this is shown below.

Prices(\$)	Equiwidth (width \$10)	Equidepth (depth=2)	Distance-based
7	[0, 10]	[7, 20]	[7, 7]
20	[11, 20]	[22, 50]	[20, 22]
22	[21, 30]	[51, 53]	[50, 53]
50	[31, 40]		
51	[41, 50]		
53	[51, 60]		

Fig: Data is partitioned by using Equiwidth & Equidepth.

Here, we examined the above table. The distance based partition is the efficient because it groups the values in such a way they are closer & also closer to specified interval.

Ex. [20, 22]

If u consider, Equidepth partition, it groups 2 distinct values i.e., [22, 50]

If u consider, Equiwidth partition, it contains equal size of intervals but some of these intervals doesn't contain any values. i.e., [31, 40]

∴ Distance based partition is the best one because this distance-based association contains the drawback i.e.,

attribute.

For ex, consider the association rule,
item-type (x, "electronic") \wedge manufacturer (x, "Apple")

\Rightarrow price (x, \$200)

Here, 'x' is the item, In reality,
we always prefer 'Apple' electronic items are
approximately '\$200' rather than 'Exactly '\$200'

\therefore To define approximate values of an attribute
& also derive distance based association rules,
we use 2 phase algorithm.

Phase-I:

Here we define interval for each cluster.

Phase-II

Each cluster is verified to identify frequent set.
Using this frequent set, we derive association
rules.

From Association Mining to Correlation Analysis:

Even the strong association rules also uninteresting
& gives the wrong information. Therefore we
require additional measures that contains
Statistical analysis.

\therefore we go for the Correlation Analysis.

Strong Association Rules are not interesting:

Consider the "All Electronics" database. we
analyse 10,000 transaction. In those 10,000

transactions, 10000

Purchase the Computer games, 7500 are Videos and 4000 transactions contains the both.

Consider Support Count is 30% & Confidence is 60%.

$\text{buys}(x, \text{"Computer games"}) \Rightarrow \text{buys}(x, \text{"Videos"})$

[Support = 40%, Confidence = 66%]

$$\text{Support}(A \Rightarrow B) = \frac{\# \text{ - tuples - Containing - A \& - B}}{\text{Total - \# - of - tuples}}$$

$$\begin{aligned} \text{Confidence}(A \Rightarrow B) &= \frac{\# \text{ - tuples - Containing - A - \& - B}}{\# \text{ - of - tuples - Containing}} \\ &= \frac{4000}{6000} = \frac{2}{3} = 0.66 = 66\% \end{aligned}$$

\therefore Rule 1 is strong association rule, but the Probability of purchasing videos is

$$\frac{7500}{10,000} = 0.75 = 75\%$$

\therefore It is greater than 66%.

So, Computer games & Videos are inversely associated - i.e., customer purchase the Computer games that decreases the Videos.

From Association Analysis to Correlation Analysis.

The above association rule $A \Rightarrow B$ is uninteresting & it gives the wrong information. So we, for the Correlation Analysis.

Correlation may be represented as

$$\text{Corr}_{A, B} = \frac{P(A \cap B)}{P(A)P(B)} \quad \text{--- (2)}$$

If the result is > 1 , then 'A' & 'B' are +ve correlated. i.e., if we increase 'A', 'B' also increases.

If the result is less than '1', then 'A' & 'B' are -vely correlated i.e., if we increase 'A' that decreases 'B'.

If the result is equal to '1' then 'A' & 'B' are independent. i.e., in between 'A' & 'B' there is no correlation.

Consider the above ex, then

Computer games = 6000

Videos = 7500

Computer games & videos = 4000

The probability of purchasing 'Computer Games' is $P(\{ \text{games} \}) = \frac{6000}{10000} = 0.60$

This is probability of purchasing 'videos' is $P(\{ \text{videos} \}) = \frac{7500}{10000} = 0.75$

The probability of purchasing both 'Computer Games' and 'videos' is

$$P(\{ \text{games, videos} \}) = \frac{4000}{10000} = 0.40$$

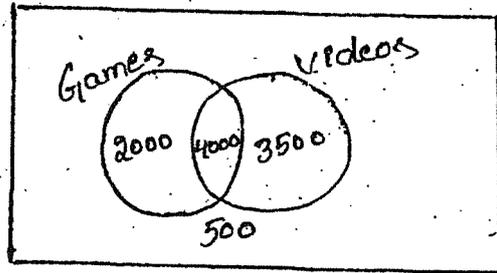
\therefore Then Correlation between Computer games &

Videos is

$$\begin{aligned} \text{Corr games, Videos} &= \frac{P(\{ \text{games, videos} \})}{P(\{ \text{games} \}) \times P(\{ \text{videos} \})} \\ &= \frac{0.40}{0.60 \times 0.75} \\ &= 0.89 \end{aligned}$$

∴ Therefore the Computer games & videos are
-vely correlated.

The transactions in DB are summarized by
Contingency Table. This is shown below.



	Games	$\overline{\text{Games}}$	Total
Videos	4000	3500	7500
$\overline{\text{Videos}}$	2000	500	2500
Total	6000	4000	10000

Fig: 2x2 Contingency Table, Summarizing Transactions Purchased.

Here ' $\overline{\text{games}}$ ' represents don't contain computer
games
 $\overline{\text{Videos}}$ represents don't contain computer video

Constraint Based Association Mining:-

Many of the DM association rules are uncovered by DM system & also discovered rules are unimportant to the end user. Therefore we go for the Constraint based association Mining.

In Constraint based association Mining, we mine association rules based on constraints. These constraints are specified by end user.

The constraints are.

i) Knowledge Constraints:-

These constraints specify the kind of knowledge to be mined.

ii) Data Constraints:-

The constraints specify the task relevant data sets.

iii) Level Constraints:-

These constraints specify the total No. of levels in concept hierarchy.

iv) Threshold Constraints:-

These constraints specify the support & confidence values.

v) Rule Constraints:-

These constraints specify what type of association rules are to be mined & also specify the no. of predicates in each rule.

Meta Rule for Mining Association Rules:-

Many of the rules are uninteresting & the wrong information. But in Meta rule, allows the user syntactic representation of association rules.

Sometimes these meta rules may contain the constraints to improve the efficiency of DM system.

The meta rule looks like the following.

$$P_1(x, y) \wedge P_2(x, z) \Rightarrow \text{buys}(x, \text{"Educational S/W"}) -$$

Here, P_1 & P_2 are predicates

$x \rightarrow$ Customer

$y, z \rightarrow$ Numeric values related to

Predicates P_1 & P_2 .

Whenever we issue this meta rule, the DM system searches for the rule that is similar to specified meta rule. Then the DM system gets the similarity & return the following association rule.

$$\text{age}(x, \text{"30 -- 39"}) \wedge \text{income}(x, \text{"30k -- 39k"}) \Rightarrow \text{buys}(x, \text{"Educational S/W"}) - \textcircled{2}$$

This meta rule is mainly used for guiding the DM process.

For ex we want to mine inter-dimensional association rules. Then the meta rule is specified as

$$P_1 \wedge P_2 \wedge \dots \wedge P_n \Rightarrow Q_1 \wedge Q_2 \wedge \dots \wedge Q_n$$

Here in this meta rule, we are using total

Predicates $P = R_1 \wedge R_2$

& also P_i where $i = 1 \dots n$ &
 Q_j where $j = 1 \dots n$

Additional Constraint Rules:

This additional Constraint rules contains set of relations. These relations are specified by the user by using aggregate functions. These rules are mainly used in multidimensional association rule Mining.

Consider the "All Electronics" multidimensional Sales db. This contains the following relations:

Sales(Cust-Name, item-Name, trans-id)

lives(Cust-Name, City, Street)

Item(item-Name, price)

transaction(trans-id, day, month, year)

Here lives, item & transaction tables are dimensional tables & fact table is Sales. Here we specify the relation by using 3.

i.e., Cust-Name, item-Name, trans-id.

Ex. we want to mine association query i.e., find the Sales of Cheap items (Sum of Price < \$100) & also find Sales of Expensive items (min of Price \$500) for Vancouver

Customer in 1999.

This query is specified in DMQL as
mines association as lives (c, "Vancouver")
from Sales

where S. year = 1999 and T. year = 1999 group
by c.

having sum(I. price) and min(I. price) will
Support threshold = 1%.

with Confidence threshold = 50%.

7. Classification & Prediction

7.1 what is classification & what is prediction:

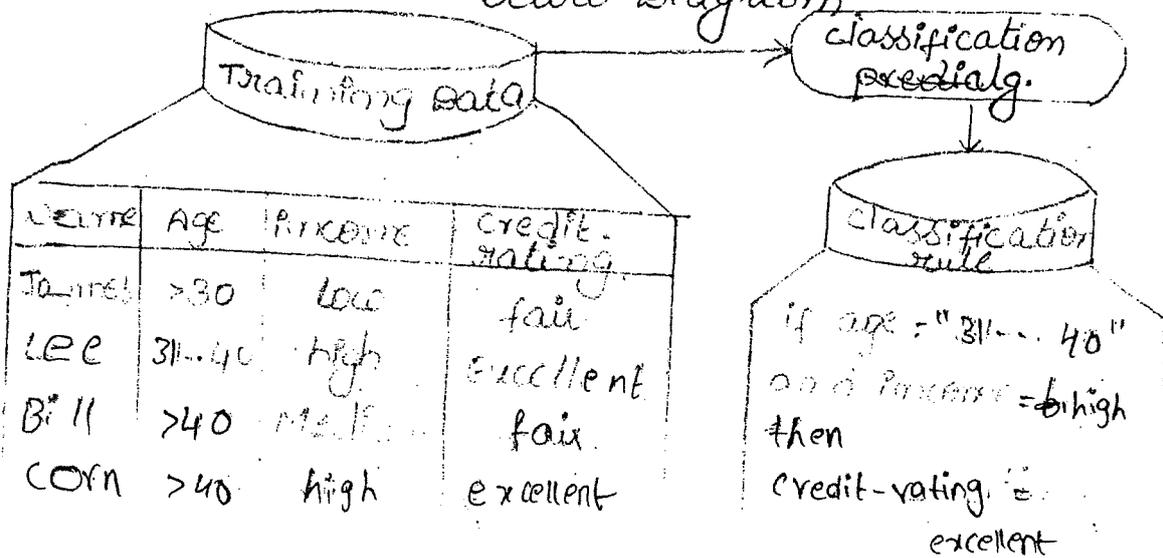
The data " " contains 2 steps.

i) construct the model by using predefined dataset: This model uses the tuples in the db & also each tuple must fall in class. This is determined by using attribute.

This attribute is called as class label attribute. Then the training data is collected & analyzed by using classification algorithm. i.e., learned knowledge is represented into classification rule.

ii) classification:

In this, classification, first verify the accuracy of the " " rule, its accuracy is ok. Then it will be applied for new data. This is shown in below diagram.



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hm.
to

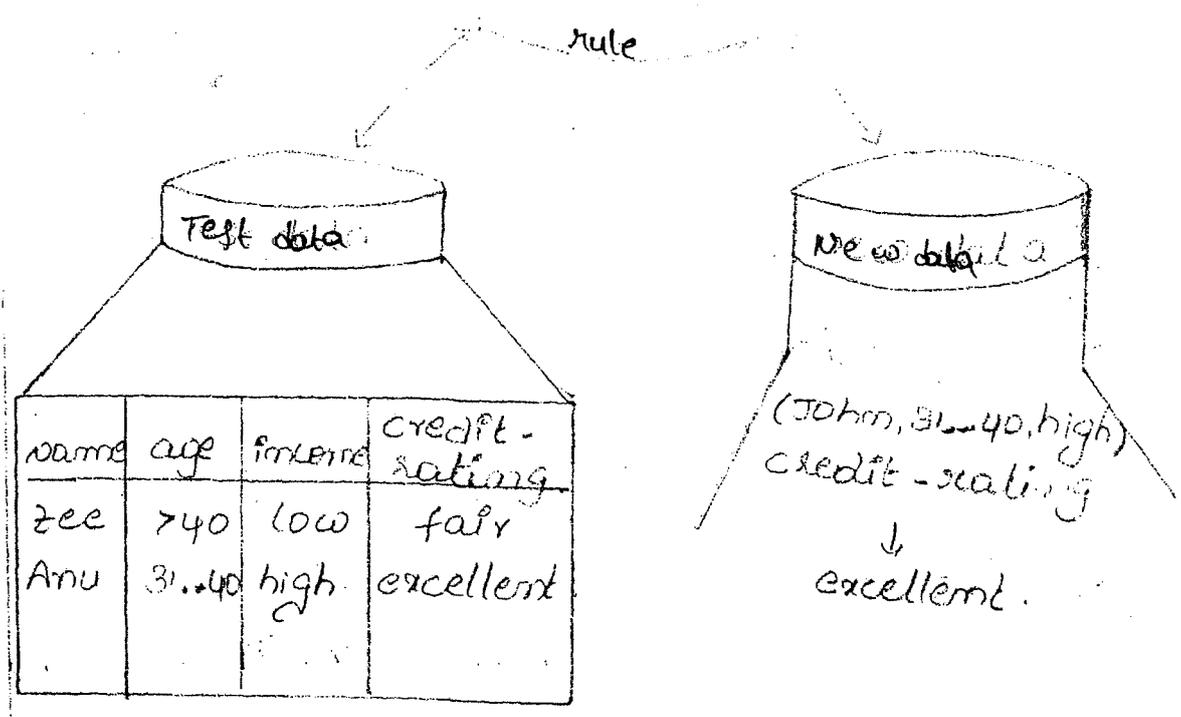


fig: 71 The Data classification.

a) Learning:

Training data is analyzed by using classification algorithm & here class label attribute is credit-rating. Then learning data is transformed into classification rule.

b) classification:

The classification rule is applied on test data, to find its accuracy. If it is accurate i.e., applied for new data.

In fig 71 (a) it analyzes the existing data of customers & using this class label of new customers is predicted.

The prediction is different from classification. The prediction allows construct the model & use the model to find the class label of unlabelled sample. The prediction uses the classification & regression.

1. Classification:

using classification we predict the discrete values.

2. Regression:

using this, we predict continuous values.

7.2 Issues Regarding classification & Predictions

we use the following issues.

7.2.1. Preparing the data for classification & prediction:

Here, we initially examine the data to improve the efficiency, accuracy, scalability of classification & prediction.

i) Data cleaning:

Here noisy or missing data values are smoothed by using several smoothing techniques.

ii) Relevant Analysis:

Here, irrelevant attributes are removed from the data before applying classification & prediction.

iii) Data Transformation:

We already know higher levels contains the more frequent items rather than the low level. This is achieved by using concept hierarchy.

The data transformation specifies generalization levels to concept hierarchy.

2319 7.2.2 Comparing Classification & Prediction Methods:

The classification & prediction methods are compared by using following methods.

1. Predictive Accuracy:

This specifies how we use the model to predict class label of new data. As well as it also, specifies the prediction must be accurate.

2. Speed:

It specifies, to predict the class label it must take the less no. of comparisons.

3. Robustness:

This specifies that we predict the class label, if the model contains some noisy data or missing values.

4. Scalability:

It specifies model is constructed from large db.

5. Interpretability:

It specifies level of understanding &

also specifies how we interpret the model.

7.3. Classification by Decision Tree Induction:

Decision Tree means flow chart like tree structure. The internal nodes specifies test on attribute. The

The branch is the outcome of the test result. The leaf node represents class. The top node of the tree is considered as a root node. The decision tree for concept hierarchy buys-computer is shown below.

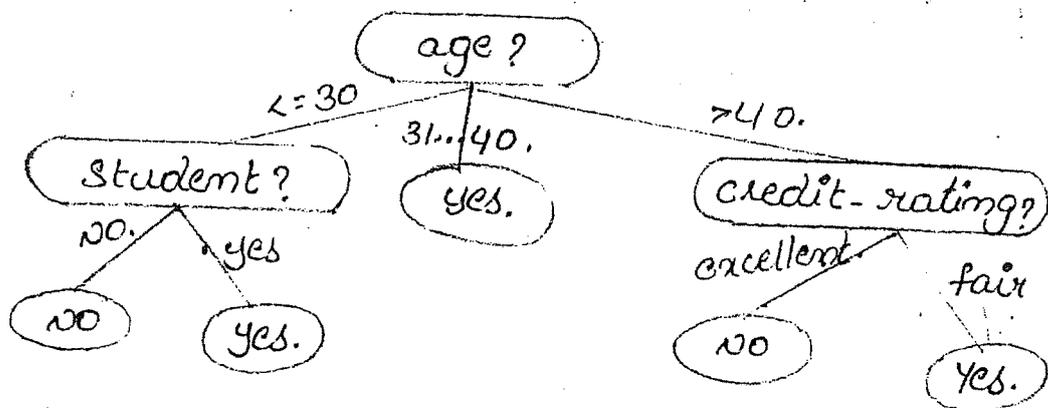


fig: 7.3. Decision Tree for Concept buys-computer indicates whether the customer buys computer or not.

The internal nodes represents test on attribute. The leaf node class (buys-computer = no or buys-computer = yes).

This decision tree contains the adv. of we classify unknown sample & also this decision tree directly converted into association

i.e. basic decision tree alg

Algorithm: generate - decision - tree.

Input: The training samples.

Output: A decision tree.

Method:

1. create a node 'v'.
2. if all the samples are same class 'c' then
3. return 'v' as a leaf labelled with class 'c'.
4. if attribute - list is empty
5. return v as a leaf labelled with most common class.
6. select test-attribute from attribute-list with highest inf. gain.
7. Label node 'v'
8. Each known value a_i of test-attribute
9. grow branch from node 'v'.
10. Let S_i be the samples in test-attribute $= a_i$.
11. if S_i is empty then
12. attach the leaf labelled with most common class.
13. else attach node & recall generate - decision - tree.

7.3.1
Basic
Decision
Tree Induction
Algorithm

7.3.1. Decision Tree Induction:

The above decision tree alg. follows the top down approach & divide & conquer approach.

The alg. is explained by using the following steps

- create a model (step 1)
- if the samples are same class, then return all the class labels related to that class (step 2,3)
- otherwise select the test-attribute by using entropy based inf. gain.
- For each known value of test-attribute, create branch
- if the samples of test-attribute are empty, then return most common class labels.
otherwise again recall the decision tree alg.

Attribute Selection:

Let 'S' be the training samples. It contains the 'm'-classes. Let 's_i' be any sample in 'S' with class label 'c_i' for i = 1...m. then expected inf. is needed to classify the given sample

$$I(S_1, \dots, S_m) = - \sum_{i=1}^m P_i \log_2(P_i)$$

Here 'P_i' is the probability of given sample fall in class label 'c_i'. It is calculated by $\frac{S_i}{S}$.

consider an attribute 'A' with values $\{a_1, a_2, \dots, a_v\}$. This attrib. 'A' partition the sample 'S' into subsets $\{S_1, S_2, \dots, S_v\}$

Let 'S_j' contains 'S_{ij}' samples of class 'c_i'. then entropy of 'A'

$$E(A) = \sum_{j=1}^s \frac{S_{ij} + \dots + S_{mj}}{S} \cdot I(S_{ij}, \dots, S_{mj})$$

∴ The inf. gain through this partition 'A' as

$$G(A) = I(S_1, \dots, S_m) - E(A).$$

The attribute with the highest inf. gain is treated as the most relevant attrib. that is used to construct the decision tree.

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Ex. consider the data tuples from 'All Electronics' database.

Page 6

Table: 7.3. Data tuples from ALL Electronics DataBase.

R I D.	age	income	Student	credit- rating.	class: buys- computer.
1.	<=30	low	no	fair	no.
2.	<=30	high	no	excellent	no
3.	31...40	low	no	fair	yes.
4.	>40	medium	yes	excellent	no.
5.	>40	high	no	fair	yes.
6.	>40	low	no	fair	yes.
7.	31..40	high	yes	fair	yes
8.	31..40	low	no	excellent	yes
9.	31..40	medium	yes	fair	yes.
10.	<=30	"	yes	fair	yes.
11.	<=30	low	"	excellent	yes.
12.	<=30	high	no	fair	no
13.	>40	low	no	excellent	no
14.	>40	high	no	fair	yes.

here class label attribute is 'buys - computer'.
 It contains 2 values {yes, no}. The class 'c₁' is
 represented by samples of 'yes', the class 'c₂' is
 " " " " 'no'.

The above table contains '9' samples of
 'yes' & '5' samples of 'no'. Therefore, the expected
 inf. is needed to classify the given sample.

$$I(S_1, S_2) = -\frac{9}{14} \log_2 \frac{9}{14} - \frac{5}{14} \log_2 \frac{5}{14}$$

$$= 0.94.$$

Then we have to calculate the entropy value
 for each attribute starting with attribute 'age'.
 This attribute contains the different branches &
 fall in 2 classes {yes or no}.

for age = "<= 30"

$$S_{11} = 2, S_{21} = 3.$$

⇒ total = 5

$$I(S_{11}, S_{21}) = -\frac{2}{5} \log_2 \frac{2}{5} - \frac{3}{5} \log_2 \frac{3}{5}$$

$$= 0.97$$

for age = "31 - 40"

$$S_{12} = 4, S_{22} = 0.$$

⇒ total = 4

$$I(S_{12}, S_{22}) = 0.$$

for age = "> 40"

$$S_{13} = 3, S_{23} = 2.$$

total = 5

$$I(S_{13}, S_{23}) = -\frac{3}{5} \log_2 \frac{3}{5} - \frac{2}{5} \log_2 \frac{2}{5}$$

$$= 0.97$$

∴ entropy of age is

$$\begin{aligned} E(\text{age}) &= \frac{5}{14} I(S_{11}, S_{21}) + \frac{4}{14} I(S_{12}, S_{22}) + \frac{5}{14} I(S_{13}, S_{23}) \\ &= \frac{5}{14} (0.97) + \frac{4}{14} (0) + \frac{5}{14} (0.97) \\ &= \frac{10}{14} (0.97) \\ &= \frac{9.7}{14} \\ &= 0.69 \end{aligned}$$

$$\begin{aligned} \text{Gain}(\text{age}) &= I(S_{11}, S_{21}) - E(\text{age}) \\ &= 0.94 - 0.69 \\ &= 0.25 \end{aligned}$$

Similarly, we can calculate the inf. gain for income is

$$\text{gain}(\text{income}) = 0.02$$

$$\text{gain}(\text{student}) = 0.15$$

$$\text{gain}(\text{credit-rating}) = 0.04$$

$$\rightarrow \text{gain}(\text{class: buys-computer})$$

Therefore, age contains the highest inf. gain. It is the most relevant attribute. Therefore, it is selected as the test attribute. Using this we construct the decision tree.

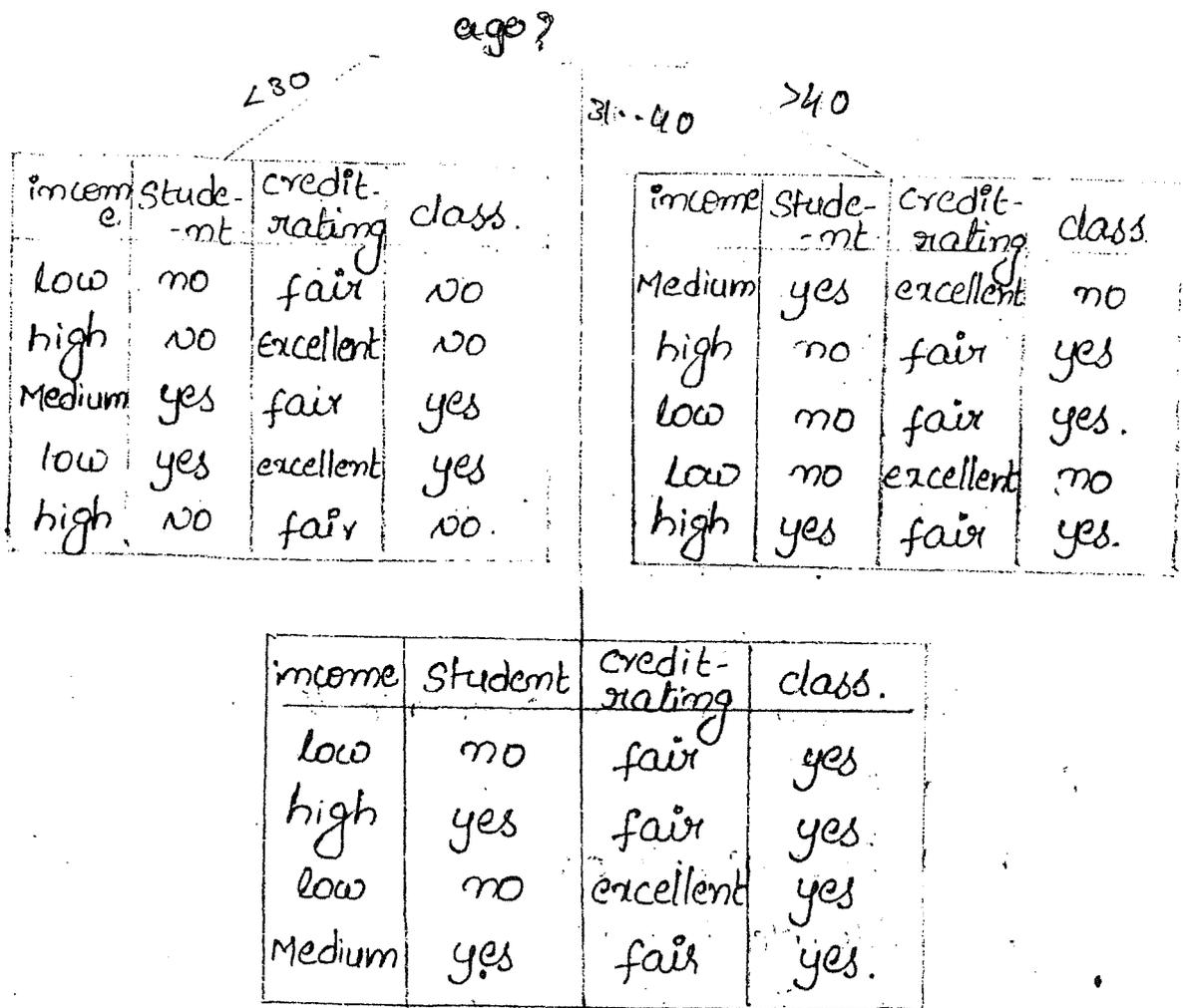
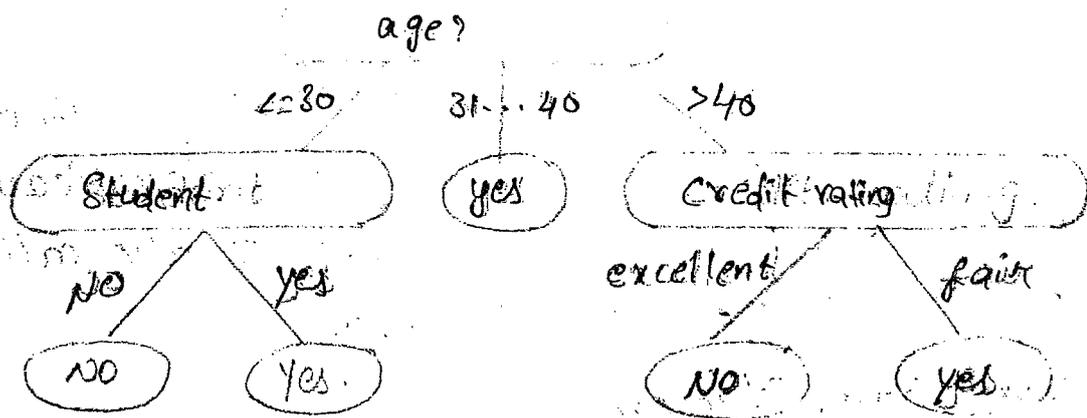


fig. 7.3. The 'age' attribute is the highest imp. gain.
 ∴ it is the root node & branches grow with 'age' values.

Here the age 31-40 contains same samples i.e., class 's'.

The final decision tree based on the alg. is shown below.

fig: Decision tree for concept buys-computer.



7.3.2 Extracting classification rules from decision trees:

The knowledge represented in decision tree extracted & represented in the form of classification IF-THEN rule. One rule is derived for one path starting from root node to leaf node

Consider the decision tree fig. 7.3.

The classification rules are,

IF age = " ≤ 30 " and Student = "NO" THEN
buys-computer = "NO"

IF age = " ≤ 30 " and Student = "yes" THEN
buys-computer = "yes"

IF age = "31..40" THEN buys-computer = "yes"

IF age = " > 40 " and credit-rating = "excellent" THEN buys-computer = "NO"

IF age = " > 40 " and credit-rating = "fair" THEN buys-computer = "yes"

7.4. Bayesian classification.

It is used to predict the class label of unknown sample. It is based on Bayes Theorem & then we apply simple Bayesian classification. The simple Bayesian classification is also called as naive Bayesian classification.

7.4.1. Bayes Theorem:

Let 'x' be an unknown sample & let 'H' be the hypothesis. Using this 'H', we predict the class label of 'x'. Finally we determine $P(H/x)$. This is called as posterior probability, i.e., probability of 'H' on condition 'x'.

For ex, consider the data samples as 'fruits'. These are described by using 'colour' & 'shape'. Let 'x' is the red colour & round shape. Then 'H' gives, x is apple. Then

$P(H/x)$ gives the confidence of x is Apple because its colour is red & it is round shape.

To determine $P(H/x)$ it requires the additional inf. i.e., $P(H)$ gives the prior probability of 'H', $P(x)$ gives the prior probability of 'x' and $P(x/H)$, i.e., $P(x)$ on condition 'H'.

Therefore using this Bayes Theorem, we define the

$$P(H/x) = \frac{P(x/H) \cdot P(H)}{P(x)} \quad \text{--- (1)}$$

7.4.2 Naive Bayesian classification:

or simple

Bayesian classification contains the following steps.

Step 1:

Consider the ' m ' class i.e., c_1, c_2, \dots, c_m . Let ' x ' be the unknown sample. Then according to naive Bayesian classification, the unknown sample ' x ' is assigned to class ' c_i ' if & only if $P(c_i|x) > P(c_j|x)$ for $1 \leq j \leq m; j \neq i$.

Thus we maximize probability of $P(c_i|x)$

But according to Bayes Theorem,

$$P(c_i|x) = \frac{P(x|c_i) P(c_i)}{P(x)} \quad (2)$$

Step 2:

The $P(x)$ is same for all the classes.

Then we maximize $P(x|c_i) \cdot P(c_i)$.

For ex, prior probabilities are not known then all the classes are equal i.e.,

$$P(c_1) = P(c_2) = \dots = P(c_m)$$

Then we maximize $P(x|c_i)$ otherwise, we maximize $P(x|c_i) P(c_i)$.

$$P(c_i) = S_i/S$$

S_i = class c_i

S = Total no. of samples.

Step 3:

If there is no dependent relationship among the attributes. Then

$$P(x|c_i) = \prod_{k=1}^n P(x_k|c_i)$$

Here, we find the $P(x_1|c_i), P(x_2|c_i), \dots, P(x_n|c_i)$

a) If A_k is categorical, then

$$P(x_k|c_i) = \frac{S_{ik}}{S_i}$$

where S_{ik} is the training samples of class ' c_i ' having value x_k for A_k .

b) If A_k is continuous-valued, then we use the Gaussian distribution.

$$\therefore P(x_k|c_i) = g(x_k, \mu_{c_i}, \sigma_{c_i})$$

$$= \frac{1}{\sqrt{2\pi} \sigma_{c_i}} e^{-\frac{(x_k - \mu_{c_i})^2}{2 \sigma_{c_i}^2}}$$

where μ_{c_i}, σ_{c_i} are the mean & standard deviation of A_k .

Step 4:

The unknown sample ' x ' is classified by determining $P(x|c_i), P(c_i)$ for each class ' c_i '.

We assign unknown sample ' x ' to class ' c_i ' if & only if $P(x|c_i) \cdot P(c_i) > P(x|c_j) \cdot P(c_j)$

for $1 \leq j \leq m, j \neq i$.

$$P(\text{credit-rating} = \text{"fair"} \mid \text{buys-computer} = \text{"yes"}) = 7/9 = 0.78.$$

$$P(\text{credit-rating} = \text{"fair"} \mid \text{buys-computer} = \text{"no"}) = 2/5 = 0.4.$$

$$\therefore P(x \mid \text{buys-computer} = \text{"yes"}) = 0.22 \times 0.22 \times 0.56 \times 0.78 = 0.02$$

$$P(x \mid \text{buys-computer} = \text{"no"}) = 0.6 \times 0.2 \times 0.2 \times 0.4 = 0.0096.$$

$$P(x \mid \text{buys-computer} = \text{"yes"}) \cdot P(\text{buys-computer} = \text{"yes"}) = 0.02 \times 0.64 = 0.0128.$$

$$P(x \mid \text{buys-computer} = \text{"no"}) \cdot P(\text{buys-computer} = \text{"no"}) = 0.0096 \times 0.36 = 0.003456.$$

Therefore,

$$P(x \mid \text{buys-computer} = \text{"yes"}) \cdot P(\text{buys-computer} = \text{"yes"}) > P(x \mid \text{buys-computer} = \text{"no"}) \cdot P(\text{buys-computer} = \text{"no"})$$

i.e., $c_1 > c_2$.

Then, unknown sample 'x' is assigned to buys-computer = "yes".

each pair of adjacent layers, it contains the weighted connection.

For ea. w_{ij} is the weight from i to j .

The multilayer means, it contains only one hidden layer. Feed-Forward means, the weighted connections never come back.

The o/p's of 1/p layer are the I/p's of hidden layer. & so on.

7.5.2 Defining a n/w Topology:

Here end user has to define the n/w topology. by specifying no. of units in i/p layer, no. of units in hidden layer, no. of units in o/p layer.

7.5.3 Back Propagation:

In this, the samples are repeated continuously. For each data sample, we modify the weights to minimize the error b/w n/w prediction & well-known class label.

The back propagation contains the following steps.

i. Initialize the weights:

Here the weighted connection b/w each pair of adjacent layers, we assign small random value.

for ex. this value ranging from
 (-1 to 1 (or) -0.5 to 0.5)

ii) Forward the selected inputs:

In this step, we find net ip & o/p for each unit & also each layer.

For ex. unit is 'j' in hidden or o/p layer. then its net ip is calculated by using a formula.

$$I_j = \sum_i w_{ij} O_i + \theta_j \quad \text{--- (1)}$$

Here w_{ij} - is weight from i to j

O_i - o/p of unit 'i'

θ_j - step of Bias. using this we differentiate the diff. ip units.

I_j - net ip for unit 'j'.

These ip's are used by hidden or o/p layers.

Then these layers apply activation function.

This is shown in below.

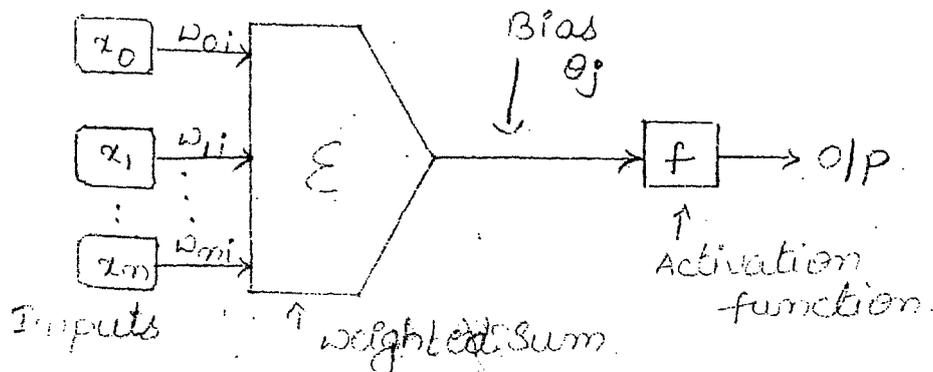


fig. 7.5.1 Activation function

Let unit 'j' & its net IIP I_j . Then o_j is the o/p for unit 'j'.

$$o_j = \frac{1}{1 + e^{-I_j}} \quad \text{--- (2)}$$

7/10 iii) Propagate error ^{Backward} forward:

To propagate the error ^{Backward} forward, we update modify the weights & Bias values.

For unit 'j' the error E_{rj} for o/p layer is calculated by using a formula

$$E_{rj} = o_j (1 - o_j) (\tau_j - o_j) \quad \text{--- (3)}$$

Here

o_j - o/p of unit 'j'.

τ_j - o/p of true value for known class label.

The error at hidden layer for unit 'j' is calculated by using the formula,

$$E_{rj} = o_j (1 - o_j) \sum_k E_{rk} w_{jk} \quad \text{--- (4)}$$

The weights are updated by using the formula.

$$\Delta w_{ij} = (l) E_{rj} o_j$$

$$w_{ij} = w_{ij} + \Delta w_{ij} \quad \text{--- (5)}$$

Here Δw_{ij} - change of weight w_{ij} .

l - linear rate. it is a constant value & ranging b/w 0.0 to 1.0.

By Bias are updated by using the formula

$$\Delta \theta_j = (-1) E_{ij}$$

$$\theta_j = \theta_j + \Delta \theta_j \quad \text{--- (6)}$$

Ex: consider the following multilayer feed forward neural n/w

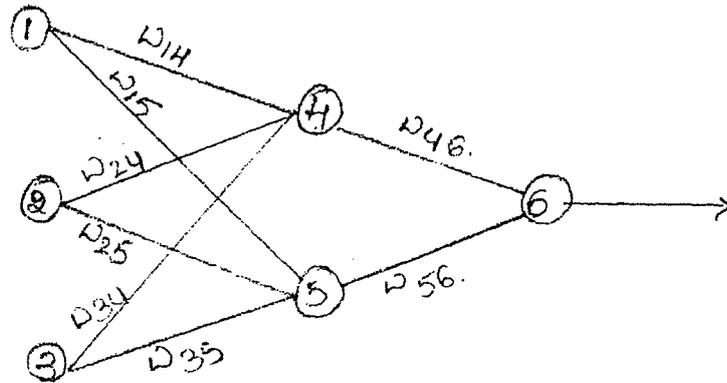


fig. 7.6. Multilayer Feed-Forward neural n/w.

The training sample,

$x = (1, 0, 1)$ whose class label is '1' & linear rate is 0.9. The initial i/p's, weights & Bias values are as shown in below.

x_1	x_2	x_3	W_{14}	W_{15}	W_{24}	W_{25}	W_{34}	W_{35}	W_{46}	W_{56}	θ_4	θ_5	θ_6
1	0	1	0.2	-0.3	0.4	0.1	-0.5	0.2	-0.3	-0.2	-0.4	0.2	0.1

Table: 7.7. Initial I/p weight & Bias values.

calculation of net I/p & o/p values:

unit j	net Input, I_j	Output, O_j
4	$0.2 + 0 - 0.5 - 0.4 = -0.7$	$1 / (1 + e^{0.7}) = 0.332$
5	$-0.3 + 0 + 0.2 + 0.2 = 0.1$	$1 / (1 + e^{-0.1}) = 0.525$
6	$(-0.3)(0.332) - (0.2)(0.525) = -0.105$	$1 / (1 + e^{0.105}) = 0.474$

calculation of Error:

unit j	Error.
6	$(0.474)(1-0.474)(1-0.474) = 0.1311$
5	$(0.525)(1-0.525)(0.1311)(-0.2) = -0.0065$
4	$(0.332)(1-0.332)(0.1311)(-0.3) = -0.0087$

calculation of weights & Bias updates:

weight/bias	new value.
w_{46}	$(-0.3) + (0.9)(0.1311)(0.332) = -0.260$
w_{56}	$(-0.2) + (0.9)(0.1311)(0.525) = -0.138$
w_{14}	$0.2 + (0.9)(-0.0087)(1) = 0.192$
w_{15}	$-0.3 + (0.9)(-0.0065)(1) = -0.3058$
w_{24}	$0.4 + (0.9)(-0.0087)(0) = 0.4$
w_{25}	$0.1 + (0.9)(-0.0086)(0) = 0.1$
w_{34}	$-0.5 + (0.9)(-0.0087)(1) = -0.5078$
w_{35}	$0.2 + (0.9)(-0.0065)(1) = 0.194$
θ_6	$0.1 + (0.9)(0.1311) = 0.2179$
θ_5	$0.2 + (0.9)(-0.0065) = 0.194$
θ_4	$-0.4 + (0.9)(-0.0087) = -0.407$

Termination condition:

Δw_{ij} becomes the small i.e., it is less than the specified Threshold range.

8/10

8. Cluster Analysis.

8.1. What is cluster analysis?

The process of storing similar objects in one cluster & dissimilar objects in another cluster. The data clusters are shown in below.

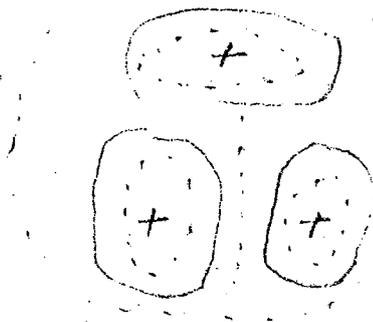


fig. 8.12-D customer data w.r to customer locations in city. Here Three data clusters are specified & centre of cluster is marked with '+'.
+
+ +

The following are the critical requirements for cluster in data mining.

① Scalability:

The clustering algorithms are efficient for low data, but DM system contains large volume of data. to handle this large volume of data, the clustering analysis alg.s must be highly scalable i.e., efficient.

8.2 Partitioning Methods.

The simplest and most fundamental version of cluster analysis is partitioning, which organizes the objects of a set into several exclusive groups or clusters.

Given a data set D , of n objects and K , the no. of clusters to form, a partitioning algorithm organizes the objects into K partitions $K \leq n$, where each partition represents a cluster.

Partitioning methods :-

- 1) K-means : A centroid-based technique.
 - 2, K-medoids : A Representative object-based technique.
- 1) K-means : A centroid-based technique :-

K-means clustering intends to partition n objects into K clusters in which each object belongs to the cluster with the nearest mean. This method produces exactly K different clusters of greatest possible distinction.

The best number of clusters K leading to the greatest separation is not known a priori and must be computed from the data. The objective of K-means clustering is to minimize total intra-cluster variance or the squared error function.

$$E = \sum_{i=1}^K \sum_{p \in C_i} \text{dist}(p, c_i)^2$$

$$E = \sum_{i=1}^K \sum_{p \in C_i} \text{dist}(p, c_i)^2,$$

where E is the sum of the squared errors.

P is the point in space representing a given object.
 c_i is the centroid of cluster C_i .

Algorithm:

K-means algorithm for partitioning, where each cluster is represented by the mean value of the objects in the cluster.

Input:

K : the number of clusters,

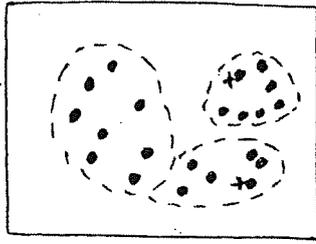
D : a data set containing n objects.

Output:

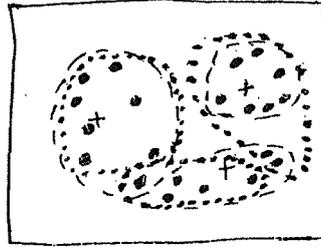
A set of K clusters.

Method:

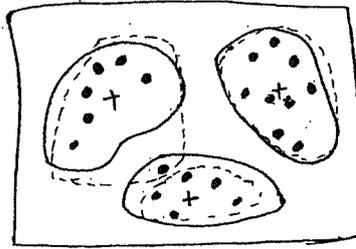
- 1) arbitrarily choose K objects from D as the initial cluster centers;
- 2) repeat
- 3) assign each object to the cluster to which the object is the most similar, based on the mean value of the objects in the cluster;
- 4) update the cluster means, i.e., calculate the mean value of the objects for each cluster;
- 5) until no change;



(a) Initial clustering.



(b) Iterate.



(c) Final clustering.

clustering of a set of objects using the k-means

Consider a set of objects located in 2-D space, as depicted in fig (a). Let $k=3$, i.e., the objects to be partitioned into 3 clusters.

According to the algorithm, we arbitrarily choose 3 objects as 3 initial cluster centers, cluster centers are marked by '+'. Each object is assigned to a cluster based on the cluster center to which it is the nearest.

Next the cluster centers are updated i.e., the mean value of each cluster is recalculated based on the centroid object in the cluster.

$$C = \sum_{i=1}^K \sum_{j=1}^n w_{ij} |p_i - o_j|$$

using this new cluster centers, the objects are re distributed to the cluster based on which cluster center is the nearest. The process of iteratively re assigning object to clusters to improve the partitioning is referred to as iterative relocation.

The time complexity of the k-means algorithm is $O(nkt)$, where n is the total no. of objects, k is the no. of clusters, & t is the no. of iterations. therefore, the method is relatively scalable & efficient in processing large data sets.

2. K-Medoids: A Representative Object-Based Technique:-

The k-means algorithm is sensitive to outliers because such objects are far away from the majority of the data, they can drastically distort the mean value of the cluster. This affects the assignment of other objects to clusters. This effect is particularly due to the use of the squared error function.

The k-medoids is an iterative clustering algorithm which iterates until each representative object is the medoid or most centrally located object of its cluster.

$$E = \sum_{i=1}^K \sum_{j=1}^n |p_i - o_j|$$

The most common realization of K-medoid clustering is Partitioning Around Medoids (PAM)

K-Medoids Algorithm (PAM) :-

Input :

K: the no. of clusters

D: a dataset containing n objects

Output :-

A set of K clusters

Method :-

- 1) Arbitrary choose K objects from D as representative objects.
- 2) Repeat
- 3) Assign each remaining object to the cluster with the nearest representative object
- 4) For each representative object o_j
- 5) Randomly select a non representative object o_{random} .
- 6) Compute the total cost S of swapping representative object o_j with o_{random} .
- 7) If $S < 0$ then replace o_j with o_{random}
- 8) Until no change.

$O(k(n-k)^2)$

The complexity of each iteration is $O(k(n-k)^2)$. For large values of n and k, such computation becomes very costly.

Advantages

Advantages:-

- * K-Medoids method is more robust than K-Means in presence of noise and outliers.

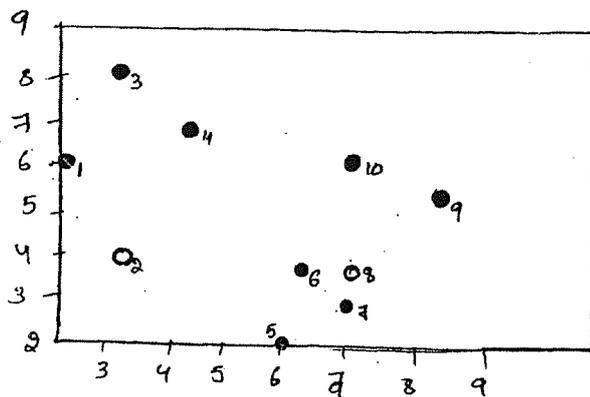
Disadvantages:-

- * K-medoids is more costly than the K-Means method
- * Like K-means, K-medoids requires the user to specify K
- * It does not scale well for large data sets.

Example:-

Data Objects

	A1	A2
O ₁	2	6
O ₂	3	4
O ₃	3	8
O ₄	4	7
O ₅	6	2
O ₆	6	4
O ₇	7	3
O ₈	7	4
O ₉	8	5
O ₁₀	7	6

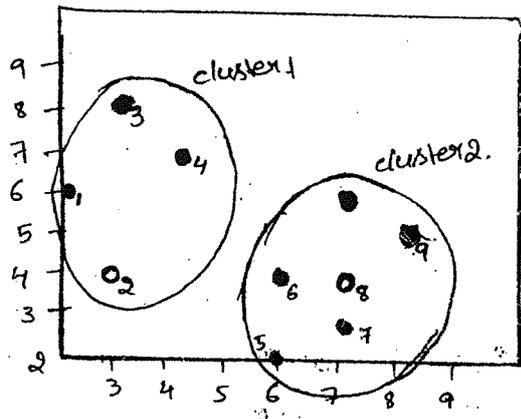


Goal: Create two clusters

choose randomly two medoids

$$O_2 = (3, 4)$$

$$O_8 = (7, 4)$$



* The absolute error/criterion (for the set of medoids (O_2, O_8))

* Assign each object to the closest representative object

* Using L1 Metric (Manhattan), we form the following clusters

$$\text{cluster 1} = \{O_1, O_2, O_3, O_4\}$$

$$\text{cluster 2} = \{O_5, O_6, O_7, O_8, O_9, O_{10}\}$$

* Compute the absolute error criterion [for the set of medoids (O_2, O_8)]

$$E = \sum_{i=1}^K \sum_{p \in C_i} |p - o_i| = |O_1 - O_2| + |O_3 - O_2| + |O_4 - O_2| + |O_5 - O_8| + |O_6 - O_8| + |O_7 - O_8| + |O_9 - O_8| + |O_{10} - O_8|.$$

* The absolute error criterion [for the set of medoids (O_2, O_8)]

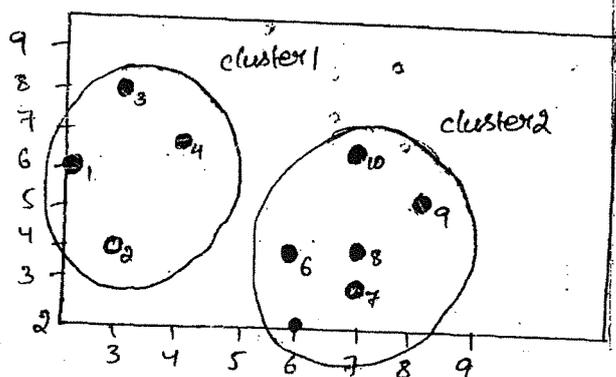
$$E = (3+4+4) + (3+1+1+2+2) = 20$$

* choose a random object O_7

* Swap O_8 and O_7

* Compute the absolute error criterion [for the set of medoids (0₂, 0₇)]

$$E = (3+4+4) + (2+2+1+3+3) = 22.$$



* Compute the cost function

$$\text{Absolute error} [0_2, 0_7] - \text{Absolute error} [0_2, 0_8]$$

$$S = 22 - 20$$

$S > 0 \Rightarrow$ it is a bad idea to replace 0₈ by 0₇

8.3 Hierarchical Methods:-

A hierarchical clustering method creates a hierarchical structure from data objects. The hierarchical clustering methods can be classified into two types. They are,

- 1) Agglomerative (bottom-up)
- 2) Divisive (top-down).

Agglomerative Hierarchical Clustering:-

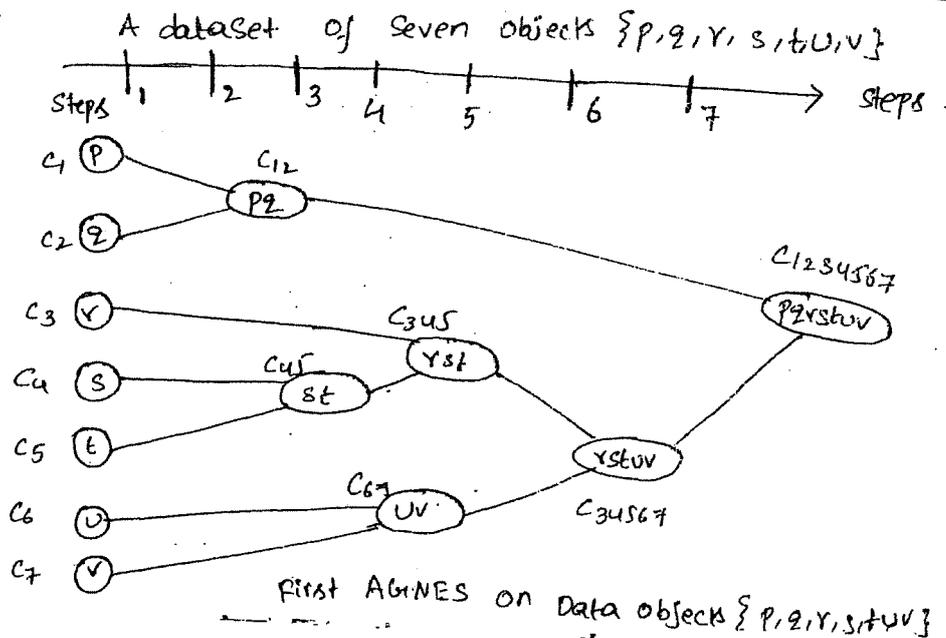
This involves merging of data objects. It initiates with each object forming its own group/cluster. For every pair of cluster, some value of dissimilarity is computed,

then the clusters are merged into larger and larger clusters until all the clusters are merged into one largest cluster or until a termination condition is reached. The merging of clusters is carried-out based on the Euclidean distance b/w any two objects from different clusters. The user can set the termination criteria by fixing the desired no. of clusters.

Divisive Hierarchical clustering:-

This clustering method is also known as top-down method. This involves division of objects cluster into smaller parts. It initiates with all the objects in the same cluster, this single cluster is splitted into smaller clusters, until each object is one cluster or until a termination condition is reached.

Example for AGNES and DIANA :- AGNES (Agglomerative Nesting)



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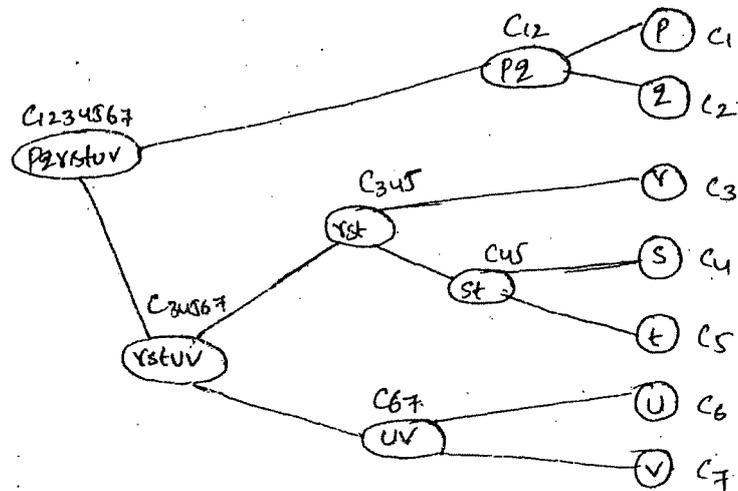
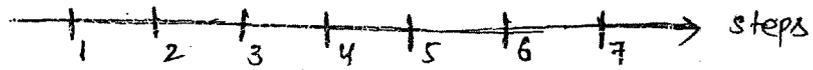
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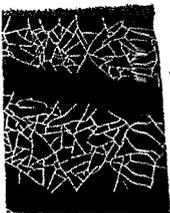
be



1, (DIANA) Divisive Analysis on Data Objects $\{P, Q, R, S, T, U, V\}$

- 1, In 'AGNES', all the objects are placed into individual clusters ($C_1 - C_7$), from the second step onwards, the objects are merged into larger objects based on the minimum Euclidean distance b/w any two objects from diff. clusters.
- 2, This process is continued until a single cluster containing all the objects is obtained.

In DIANA, all the data objects are first placed into a single cluster ($C_{1234567}$). This single cluster is then divided into smaller clusters based on maximum Euclidean distance b/w closest neighbouring object in the cluster. This process continues until each object is placed in an individual cluster.



8.3.2 Distance Measures in Algorithmic Methods :-

Whether using an agglomerative method or a divisive method, a core need is to measure the distance b/w 2 clusters, where each cluster is generally a set of objects.

Four widely used measures for distance b/w clusters are as follows.

$$\text{minimum distance: } \text{dist}_{\min}(C_i, C_j) = \min_{p \in C_i, p' \in C_j} \{ |p - p'| \}$$

$$\text{Maximum distance: } \text{dist}_{\max}(C_i, C_j) = \max_{p \in C_i, p' \in C_j} \{ |p - p'| \}$$

$$\text{mean distance: } \text{dist}_{\text{mean}}(C_i, C_j) = |m_i - m_j|$$

$$\text{Average distance: } \text{dist}_{\text{avg}}(C_i, C_j) = \frac{1}{n_i \cdot n_j} \sum_{p \in C_i, p' \in C_j} |p - p'|$$

where $|p - p'|$ is the distance b/w 2 objects or points

m_i is the mean for cluster C_i ,

n_i is the no. of objects in C_i

they are also known as linkage measures.

When an algorithm uses the minimum distance, $\text{dist}_{\min}(C_i, C_j)$ to measure the distance b/w clusters, it is called as "nearest-neighbor clustering algorithm".

An agglomerative hierarchical clustering algorithm uses the minimum distance measure is also

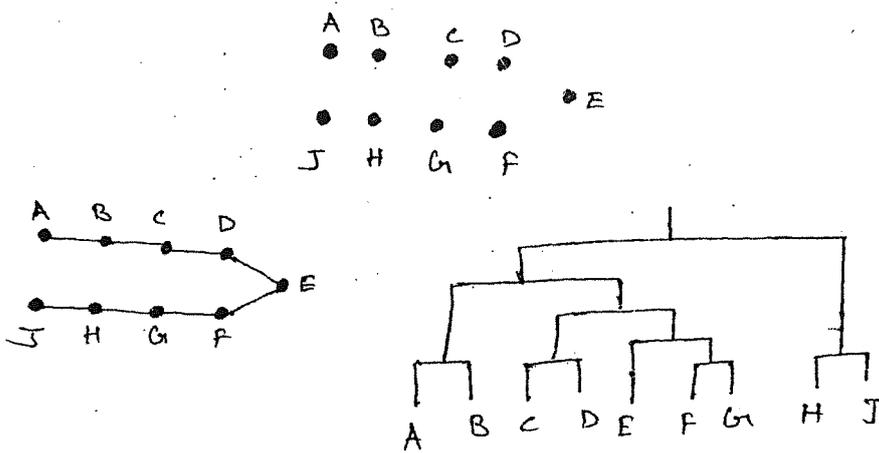
Called a "minimal spanning tree algorithm". The minimal spanning tree is the one with the least sum of edgeweights.

When an algorithm uses the maximum distance $d_{max}(c_i, c_j)$, to measure the distance b/w clusters it is called as "farther-neighbor clustering algorithm".

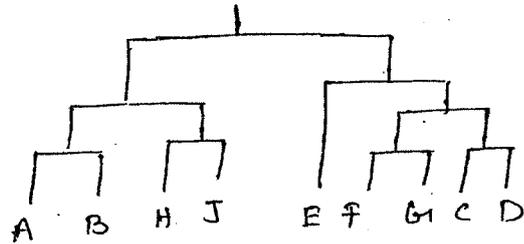
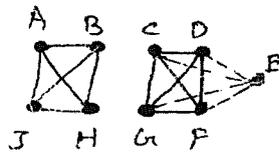
The use of mean or average distance is a compromise b/w the minimum & maximum distance and overcomes the outliers sensitivity problem.

Example :-

let us apply the hierarchical clustering to the data set $\{A, B, C, D, E, F, G, H, J\}$



clustering using single linkage



clustering using complete linkage.

8.3.3 BIRCH: Multiphase Hierarchical clustering using clustering feature trees

BIRCH (Balanced Iterative Reducing and clustering using Hierarchies) method integrates both hierarchical clustering methods, in initial stage and iterative partitioning clustering method at final stages. In other words, the o/p generated from hierarchical methods serve as I/p or preprocessing step for iterative partitioning.

Advantages of BIRCH when compared to Hierarchical Methods

- 1) BIRCH method improves the scalability of clustering.
- 2) It has the ability to undo the mistakes which have occurred in previous phases.

In BIRCH method, for an input data points, the following three factors are calculated.

1) Centroid (C)

It determines the center of the cluster. It is

Calculated using the formula, $G_c = \frac{\sum_{i=1}^n z_i}{n}$

2, Radius (r)

It determines the average distances of the I/p point from the centroid.

$$r = \sqrt{\frac{\sum_{i=1}^n (z_i - G_c)^2}{n(n-1)}}$$

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3, Diameter (D)

It determines the average pair wise distance in a cluster.

$$D = \sqrt{\frac{\sum_{i,j=1}^n (z_i - z_j)^2}{n(n-1)}}$$

The radius and diameter factor determines the level of association of the cluster around the centroid.

Clustering Feature and Clustering Feature Trees:-

In BIRCH method, clustering Feature (CF) is defined as a three-dimensional vector, which contains the info about the clusters of objects that have been discarded or compressed in a summarized form.

The 'CF' for 'n' data points spread over d-dimensions can be represented as,

$$CF = (n, L_{sum}, S_{sum})$$

where

$n \rightarrow$ no. of data points (or) Objects.

$L_{sum} \rightarrow$ Linear sum of data points,

$S_{sum} \rightarrow$ Square sum of data points.

Clustering Feature Tree (CFT) is height-balanced tree which is used to represent clustering features of individual clusters in an hierarchical fashion.

The non leaf nodes maintain the info. about the sums of the clustering features of their child nodes and thereby summarizing the clustering info. of these child nodes.

The size of a clustering feature tree is dependent on two factors.

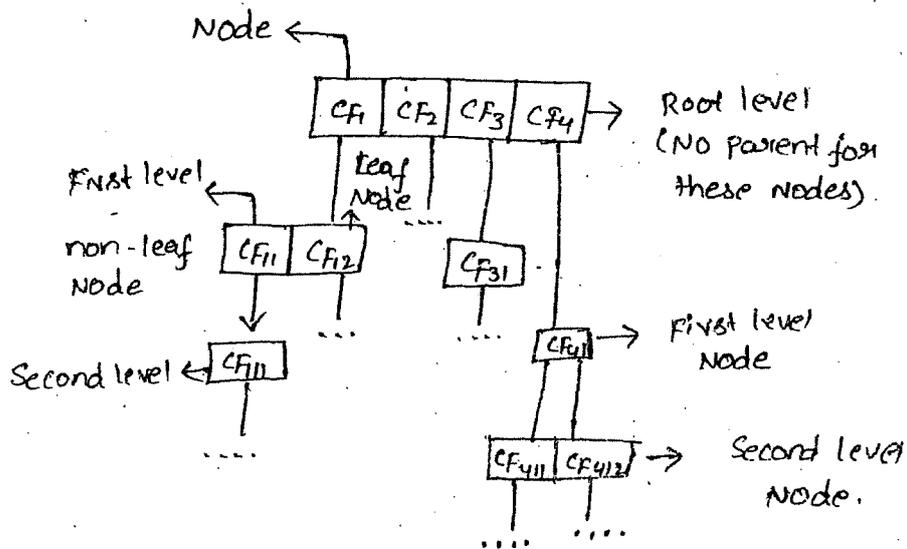
1. Branching factor (B)

This factor decides the maximum no. of child nodes for a non leaf node.

2. Threshold (T)

Threshold decides the maximum diameter that a subcluster i.e., a collection of non-leaf and its child nodes.

Example of CF Tree



BIRCH method is used to generate efficient clusters using those resources that are available. This method applies multiphase clustering technique which generates good clustering with minimum intervention from I/O and minimum amount of main memory for storage.

8.3.4 Chameleon: A Hierarchical clustering Algorithm using Dynamic Modeling :-

This method uses dynamic modeling. It basically constructs a sparse k -nearest neighbor graph, then partitions the graph into pieces and then clusters the pieces together. produces more natural clusters than DBSCAN. uses density measurements to determine the k -nearest neighbor.

Notes

adapt to the characteristics of the data set to find the natural clusters.

level
parent for
se nodes)

use a dynamic model to measure the similarity
between clusters.

1st level
node

main property is the relative closeness and relative
inter connectivity of the cluster.

Second level
node.

Two clusters are combined if the resulting cluster
satisfies certain properties with the constituent clusters.

The merging scheme preserves self-similarity

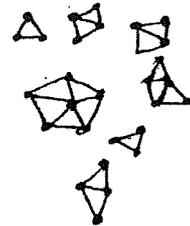
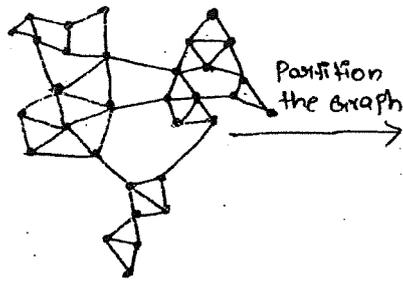
efficient clusters of the areas of application is spatial data.

table. This method

k-nearest-neighbor graph.

tree which generates
intervention from I/O
memory for storage.

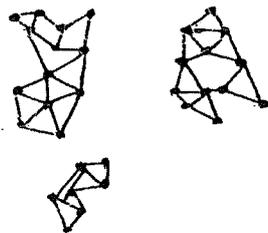
reconstruct
parent
graph



Clustering Algorithm

merge
partition

Final clusters



modeling. It basically
is a graph, then partitions
into clusters the pieces
smaller clusters than DBSCAN.

Hierarchical clustering Based on k-nearest neighbors
determine the k-nearest
neighbor modeling.

Chameleon Steps :-

Preprocessing :-

Represent the data by a graph

1, Given a set of points, construct the k -nearest neighbor (k -NN) graph to capture the relationship b/w a point and its k -nearest neighbours.

2, concept of neighborhood is captured dynamically.

Phase 1 :-

use a multilevel graph partitioning algorithm on the graph to find a large no. of clusters of well-connected vertices.

Each cluster should contain mostly points from one "true" cluster i.e, is a sub-cluster of a "real" cluster.

Phase 2 :-

use hierarchical agglomerative clustering to merge sub-clusters.

1, Two clusters are combined if the resulting cluster shares certain properties with the constituent clusters.

2, There are two key properties used to model cluster similarity.

* Relative Interconnectivity :-

Absolute interconnectivity of two clusters normalized by the internal connectivity of the clusters.

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Noise

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* Relative closeness :-

Absolute closeness of two clusters normalized by the internal closeness of the clusters.

nearest
relationship
is.

similarly

4 Density-Based Methods :-

Density based clustering methods have been developed to determine clusters with arbitrary space. clustering based on density, such as density-connected points.

them on the
well-connected

Major Features of Density-Based Methods :-

non "true"
sets.

- 1) Discover clusters of arbitrary shape.
- 2) Handle noise.
- 3) One scan.
- 4) need density parameters as termination condition.

merge

there are three different methods of density-based

cluster

clustering. they are:

clusters.

1) DBSCAN

cluster

2) OPTICS

3) DENCLUE

8.4.1 DBSCAN :-

normalized

Density-Based Spatial Clustering of Application with Noise (DBSCAN) is a density-based clustering algorithm.

In DBSCAN a cluster is a set of maximum density

connected points DBSCAN finds arbitrary-shaped clustering alg: and finds arbitrary-shaped clusters in spatial DBs with noise and expands regions having enough high density into clusters. To better understand the concept of DBSCAN method we first need to know some definitions.

Noise

Noise is defined as the set of objects in N which do not belong to any cluster.

1, ϵ -neighborhood of an object:-

The ϵ -neighborhood of an object is defined as for a given positive radius, the neighborhood of an object must be within the radius (ϵ).

2, CORE object:-

An object is said to be a core object, if the ϵ -neighborhood of an object contains at least a minimum no. of objects (Minpts).

3, Directly-Density Reachable:-

An object 'e' belonging to a set of objects, N is directly density-reachable from an object 'f', if 'f' is a core object and e is within ϵ -neighborhood of f.

4) Density-Reachable:-

Given a neighborhood ϵ and minipnts in a group of objects N and when there is a series of objects e_1, e_2, \dots, e_n where $e_1 = f$ and $e_n = e \Rightarrow e_{a+1}$ is directly density reachable from e_n with respect to ϵ and minipnts, for $1 \leq a \leq n$, $e_a \in N$ then an object e is density reachable from object f .

5) Density-Connected:-

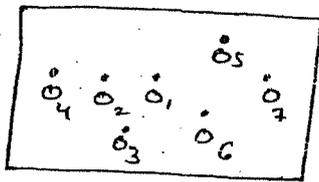
Given ϵ and minipnts in a group of objects N and when there is an object 'g' belonging to $N \Rightarrow$ both e & f are density-reachable from 'g' with respect to ϵ and minipnts. then an object 'e' is density-connected to object f .

6) Density clusters:-

A cluster with respect to ϵ and minipnts is a non-empty subset of N containing maximum no. of density connected objects with respect to density-reachability.

Example:-

let us assume that minipnts = 5 and $\epsilon = 2$ cm.



objects before clustering.

steps for finding clusters:-

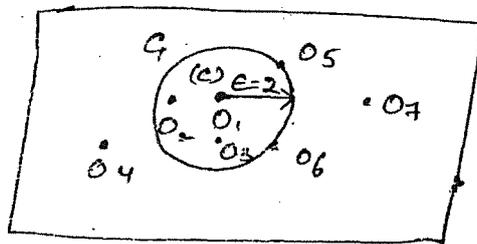
1. Select an unclassified object o_1 . The objects that are in the neighborhood of o_1 along with the radius ϵ are,

$$N_E(O_1) = \{O_1, O_2, O_3, O_5, O_6\}$$

since, $N_E(O_1) \geq \text{minpts}$ i.e., $5 \geq 5$

O_1 is considered as a core object (c)

Thus, these set of points together form a cluster, which is assigned a cluster-id (C1). This is shown in below.



O_1 Neighbouring Objects

2, Now, select the object O_2 from the set. The neighbouring objects of O_2 are

$$N_E(O_2) = \{O_1, O_2, O_3, O_4\}$$

since $N_E(O_2) < \text{minpts}$ i.e., $4 < 5$, the object O_2 is considered as a 'noise' object (N) and therefore no cluster is formed.

3, select the object O_3, O_4, O_5 from set. The neighbouring objects of O_3, O_4, O_5 are

$$N_E(O_3) = \{O_1, O_2, O_3, O_6\}, N_E(O_4) = \{O_2, O_4\}$$

$$N_E(O_5) = \{O_1, O_5, O_6, O_7\}$$

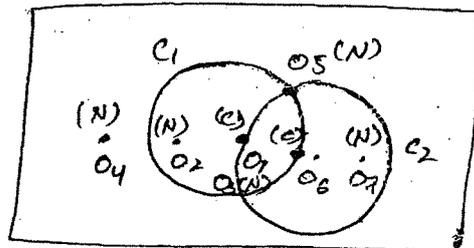
since $N_E(O_3) < \text{minpts}$, $N_E(O_4) < \text{minpts}$, $N_E(O_5) < \text{minpts}$

so, these are considered as noise objects (N), and therefore no cluster is formed.

Finally, select the object O_6 from the Set. The neighboring objects of O_6 , are

$$N_e(O_6) = \{O_1, O_3, O_5, O_6, O_7\}$$

Since, $N_e(O_6) \geq \text{minpts}$ i.e., $5 \geq 5$ the object O_6 is considered as core object. Thus, these set of points form a cluster, ~~is~~ assigned a cluster-id (C_2).



O_1 & O_6 objects with their neighbors forming clusters C_1 & C_2 .

The resultant DBSCAN clustering in which there are two clusters as,

$$C_1 = \{O_1, O_2, O_3, O_5, O_6\} \text{ \& } C_2 = \{O_5, O_6, O_7\}$$

$$C_2 = \{O_1, O_3, O_5, O_6, O_7\}$$

where O_1, O_6 are core objects and O_2, O_3, O_4, O_5, O_7 are noise objects.

SQL OPTIMIS:-

Algorithm:-

DBSCAN : a density-based clustering algorithm.

Input:

D : a dataset containing n objects,

ϵ : the radius parameter &

minpts : the neighborhood density threshold.

output:- A set of density-based clusters;

Method:-

1. mark all objects as unvisited;

do

randomly select an unvisited object p ;

mark p as visited;

if the ϵ -neighborhood of p has at least $minpts$ objects

create a new cluster C and add p to C ;

let N be the set of objects in the ϵ -neighborhood of p ;

for each point $p' \in N$

if p' is unvisited

mark p' as visited;

if the ϵ -neighborhood of p' has at least $minpts$ points

add those points to N ;

if p' is not yet a member of any cluster,
add p' to C ;

end for

output C ;

else mark p as noise;

until no object is unvisited;

Classification according to the
database mined.

Class according to the
knowledge mined.

Class according to the
techniques utilized.

Class according to the
applications adopted.