





We first discussed a hierarchical model for social networks, and then generalized it to encompass all the objects in our problem domain. In social network analysis, social networks are commonly represented by directed or ‘link’ graphs. The vertices  $i, j \in I$  represent the individuals in the network, while the relationships among them are depicted by the edges. The relationship between an individual  $i$  and  $j$  is denoted by the link from vertex  $i$  to  $j$ . It is important to quantify the association among members of the network, to achieve this, nonnegative real-valued weights are assigned to each link. The weights can be interpreted as the level of ‘endorsement’, one member grants another. For instance, the weight on the edge  $i \rightarrow j$  can be viewed as how much member  $i$  endorses or respects member  $j$ ; its value significantly indicates how  $j$  is rated among its peers in the network. This abstractive illustration showed the importance of individuals’ perception within a social media network. The aggregation of individual judgment is what bring about a finalistic rating within the network. Hence, formation of judgement graph can be achieved by processing link graph as in the case of a typical tree structure but in this case leaves are denoted with vertices. In the hierarchical model representation (see fig. 3.1) of the social network, the individuals are the leaves, while the intermediate levels are referred to as the interior nodes.

Our model easily generalized to all the objects in our problem domain. They can be represented as hierarchical social networks. For instance, in the Web, the web pages are the individuals, the hyperlinks represent the network while hierarchy represents the domains and sub-domains. This representation can be carried in diverse form depending on the subject domain. It is applicable I any knowledge representation and denotations. For instance, in a citation index, publications can be referred to as individuals, the references from one publication to another can be denoted as the network structure while the classification by subjects and sub-subjects defined an explicit hierarchical structure.

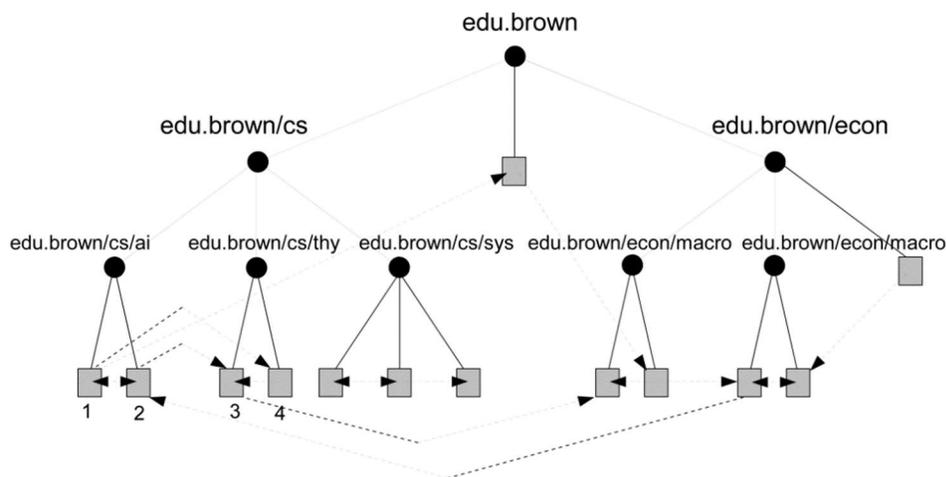
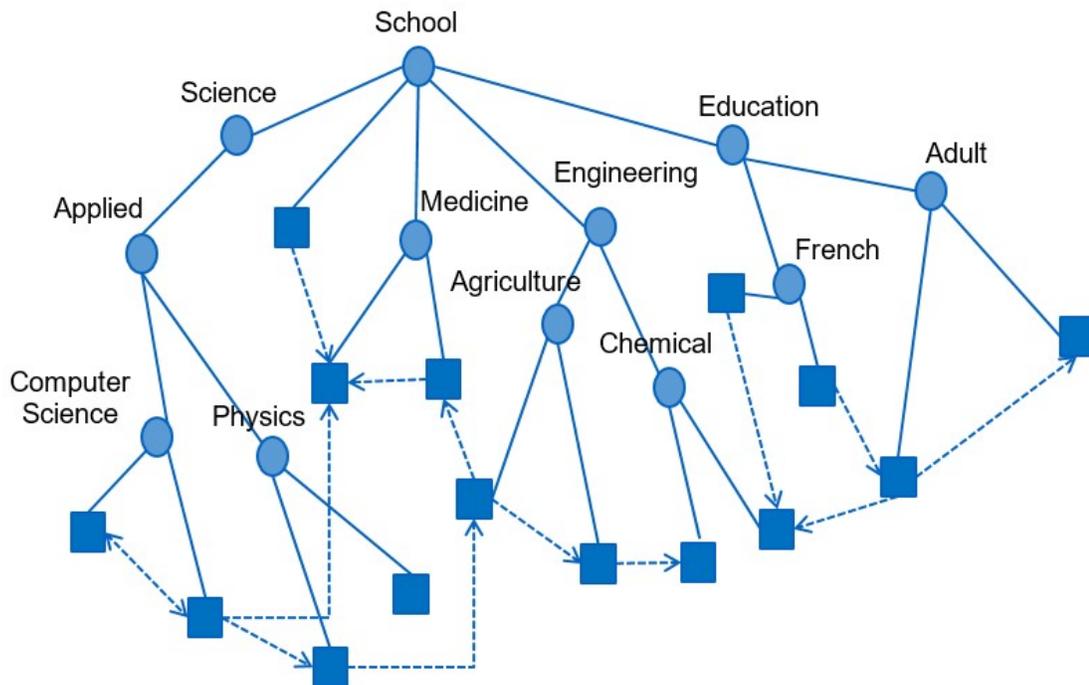


Fig..1. Hierarchical Representation of a Web domain<sup>1</sup>

<sup>1</sup> From Greenwald and Wicks (2006)

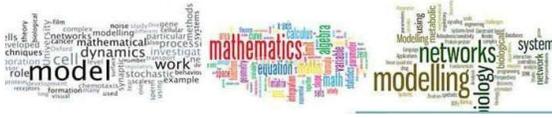
We incorporated user’s bias into the hierarchical structure by introducing user’s criteria at the leaves of the tree (solid rectangles). These criteria are the qualities identified by users as important; they form the basis for ratings and ranking. A sample personalised hierarchical model is shown in fig. 3.2—a network of the faculties and departments within a school hierarchy. At the top of the tree hierarchy are schools followed by faculties and departments to form a complete tree structure. The departments, indicated by solid circles, are the individuals in this society. The link relationship within the school hierarchy is represented by solid lines.



**Fig. .1. A Sample Hierarchical Representation of Our Proposed Model**

The relationships between criteria are indicated by dashed lines with arrowheads. Since individuals are expected to represent what they want in the best way it makes meaning to them, such a representation varies from user to user; hence, the hierarchical representation of the user domain together with the user-defined criteria could be modeled in the form of variant concept maps. Implicitly, various decision questions such as ‘is-important’, were included in the model to enable us arrived in a destination that closely matches the individual’s expectation, based on the selected criteria. We referred to this sample hierarchical model in the course of the formulation and analysis of our algorithm. We noted that the general principles of our approach can be extended to any kind of network, which can be represented by a hierarchical structure.

Modelling a hierarchical representation structure like we have in the school example above underlines the importance of every individual with the network structure. The strength of each individual cumulatively form the overall rating of each object within the hierarchy. To have a good school, all the element within school structure must be taken into consideration. In addition, the individual has varying interest while making decision regarding the all factors and elements within school ecosystem. Differential in preferences would definitely lead to different ratings and choices.



Having formulated the personalized hierarchical representation of our search object, we proceeded to develop a recursive ranking algorithm that explicitly took advantage of this structure.

## 2. ENHANCED RECURSIVE RANKING ALGORITHM

In the preceding section, we described a model for a hierarchical social network that incorporates user's bias into the structure. Here we discuss the process of designing and development of the algorithm. The algorithm is designed to exploit the underlying hierarchical topology of the network as presented in the previous section. The rating and ranking were personalized using user-defined preferences as biases. In order to ensure that our algorithm adequately solves our research problems, we proposed requirements, which it must satisfy:

1. **Hierarchical structure of the network.** The algorithm must be designed to take advantage of the inherent hierarchical structure of social network. This implies ease of transverse through the hierarchy of trees, in a Tree, Node, and Level (TNL) manner.
2. **Personalization of ranking.** The algorithm must reflect and address the issue of individual preferences by implementing a ranking procedure that incorporates users' biases within the network structure.
3. **Resistance to link-spamming.** Link-spamming is refers to exercises that are carried out with the aim of altering the ranking of a node, by creating many false nodes to that node. The false nodes are also called sybils (Cheng and Friedman 2006). A common example of link spanning is Web spamming (Gyongyi and Garcia-Molina 2004). Link-spamming presents a serious problem in ranking systems, especially in large networks. It reduces the efficiency of the ranking algorithm. A good ranking algorithm should demonstrate remarkable resistance to link-spamming.
4. **Scalability.** Problem domain that are represented in link-tree hierarchical structure are usually complex in nature. As a result, computations can be resource-intensive as the problem size increases. The algorithm must be scalable to handle the dynamic growth of the problem representation. Recursion presents the best approach. It would make the proposed solution more scalable, for efficient handling of future growth within the network. With recursion, the algorithm could be broken down into modules which are developed into components and then recursively applied across the domain.

Among the existing ranking algorithms reviewed in the previous chapter, the most similar to our proposed algorithm is QuickRank algorithm, which served as the prototype. We modified the QuickRank algorithm to incorporate user's bias. This way, we retained the desirable advantages of the algorithm, while providing adequate solution to the research problems. The selection of QuickRank was based on its desirable characteristics:

- QuickRank was designed with ease of transverse through the hierarchy in a Tree, Node and Leaf (TNL) manner. Hence, it is very suitable for developing solution to problems that can be represented in TNL fashion. It is able to take advantage of the inherent hierarchical structure of the social network resolving accordingly from leave to tree and vice versa as the case may be.
- QuickRank is recursive in nature, offering scalability, ease of parallelization, and ability to update. With this, problems can be broken down into modules with each module as a unit serving as input into another module or the main.
- QuickRank has proven to be remarkably resistant to link-spamming, and efficiently handles future growth within the social network.





Higher weight from node A to B means individual A has high level of endorsement for individual B but not vice versa. A modified form of link graph which possesses positive diagonal values is referred to as judgement graph. In such graph, individual judgement is represented by columns.

#### 2.1.4 BaseRank procedure

BaseRank is a ranking mechanism that computes and aggregates all the ranking parameters for the elements in the hierarchical structure that was described in previous section. It forms the core of our ranking algorithm by precomputing values for all the objects in our problem domain, we can then apply users' bias on the precomputed value to achieve the desirable results. Some of the BaseRank procedures that were evaluated in the course of this research are in-degree, out-degree, eigenvector centrality and PageRank (Greenwald and Wicks 2006). The output from BaseRank procedure is referred to as posterior ranking result and is a key input in the development of a recursive ranking algorithm. This output from BaseRank is then normalized into a probability distribution for easy integration. With BaseRank procedure, all the judgements of individual within a network can be aggregated and computed to generate only one aggregated posterior ranking. For illustration, assuming there is a graph  $R$  and a prior ranking  $\langle r \rangle$ , by applying Bonacich's hypothesis, an inference to a collective judgement can be expressed as  $r' = Rr$ .

#### 2.1.5 Parameterisation

This refers to a process of formulation and describing the parameters that are used with a model or system. It is a major requirement in the defining the specification and characterization of a model. The end product of parameterisation is usually in the form of expression or equation which can be applied as part of the components of the model. For instance, QuickRank is parameterised by BaseRank procedure.

#### 2.1.6 Bonacich's hypothesis

Bonacich's hypothesis deals with the how to rate relative importance of each individual in a network. It does not just consider the ratings of individuals by its peers but also the inherent rating of the peers themselves. Consequently, the importance of an individual in a network is determined by the importance of those that endorsed him/her. It further means that the strength of an individual endorsement is directly proportional to the ranking of the individuals in the network that are making the endorsement. In other words, the collective ranking of an individual in a network is the total sum of individual judgments with the weight determines by the implicit ranking. Bonacich's metric, based on this hypothesis, is also known as eigenvector centrality.

#### 2.1.7 Peer-review principle

(Greenwald and Wicks 2006) Peer-review principle states that endorsements among peers of individuals on the same sub-tree or community is more appropriate and should be taking seriously. Individuals at the same level in a network has many interaction and relationship and therefore rating in such situation should be fairer. In peer-review every other endorsement can be seen as approximation as the individuals may not possess enough information to make the right judgement while rating other individuals outside their community. As a result, while designing ranking algorithm, more weight should be assigned to ratings by individuals in the same community than others. At the heart of the QuickRank algorithm are **Bonacich's hypothesis** and "**peer-review principle**".





### 3. THE DEVELOPMENT OF ENHANCED RECURSIVE RANKING ALGORITHM

As already stated, the personalized hierarchical model of the underlying network directly incorporates user bias. ERRA, working on the model, introduces the user bias as a real-valued weight at both the leaves and nodes levels of the tree. To illustrate the algorithm, we used tree analysis which consist of judgement graph  $R$  that contains vertices which are represented as leaves of tree  $T$ . There is the assumption that the link graph is pre-computed to produce judgement graph. The ranking of the individuals in the network is then equivalent to aggregating the individual judgments into a single collective ranking. According to Greenwald and Wicks (2006), the global or resultant ranking within a social network hierarchy can best be computed by first computing the local ranking of all the nodes in the network, and then use chain rule to aggregate all the local. This follows the important fact, that all rankings are considered as a form of probability distribution before they are normalised. The ranking is done by firstly at the leaf level based on the link information that is obtained from the processing of marginal subgraphs. The local rankings are then propagated and aggregated across the network in other to derive the absolute global ranking. Nodes ranking are computed recursively, to determine a node ranking, one needs to aggregate the local ranking of the node and the local ranking of all the associated interior nodes and leaves. A BaseRank procedure is used to compute the local rank. Prior ranking of the leaves is taken as an input, and posterior distribution as an output.

Our algorithm is bottom-up, starting from the bottom of tree  $T$ , the algorithm starts by collapsing the most posterior node. Then it progress gradually by continuously identifying the next nodes and collapse accordingly. This is done progressively until there are no more nodes to collapse in which case, the tree  $T$  becomes a leave node. Assuming there is a node  $n$ , the process of collapsing this node involve three major steps as highlighted below:

- i. Computing a local ranking at  $n$  (the rankings of  $n$ 's children); this including the computation of local ranking for all the nodes that are associated with  $n$  both interior and posterior.
- ii. Computing  $n$ 's ranking and judgement; this is achieved by aggregation of all the local rankings and judgements of  $n$ 's children respectively.
- iii. Aggregating the bias for  $n$ 's children into only one weighted value which is an associate of  $n$ . This is where the ERRA differs from QuickRank—QuickRank does not include this step. User's bias  $\beta$  is a pre-computed weighted value that is applied to all  $n$ 's children. The value usually ranges from 0 to 1 ( $0 < \beta \leq 1$ , where  $\beta$  is the user's bias at leaves level). User's bias of 1 means the QuickRank value for the leaf or node stays the same. If the user's bias for all the leaves, interior and posterior nodes is equal to 1, then our Enhanced Recursive Ranking Algorithm returns the same ranking results as QuickRank. We require that the user bias be normalised (the reason and the method for bias normalisation are discussed in section 3.3).

It is easy to see that ERRA just like QuickRank is a well-defined algorithm. The global ranking does not depend on the order of computation of the local ranking. This is made possible because the algorithm adopted the approach of propagating local computations from bottom to the top of the tree while to generate resultant global values. ERRA is designed to be deterministic and terminate when there are no more nodes to collapse and the root node becomes a leaf node.



**Data Structures:** In ERRA, the two main input are;  $T_n$  which is a subtree of  $T$  at node  $n$ , and the user bias which defines user specific presence. The output is return in the form of data structure as enumerated below:

- i. Local ranking of the leaf nodes that are in the tree up to the support point  $T_n$
- ii. Computed judgement, that is the mean of all judgements of  $T_n$ 's leaves weighted by the ranking computed in (i).
- iii. A user bias, computation of the mean value for biases of  $T_n$ 's leaves.

QuickRank returns only (i) and (ii).

At the leaf node  $n$ , the ranking follows a normal probability distribution with all weights on  $n$  (denoted by  $e_n$ ), the judgment is given by  $R_n$  and bias as  $\beta_n$ .

**Computing Local Ranking:** Our algorithm relies on QuickRank's BaseRank procedure such as In-degree, Out-degree, Eigenvector centrality or PageRank (see Greenwald and Wicks 2006). BaseRank procedure takes  $n$ 's local prior ranking and a local judgment graph  $M$  as input. For  $j$  and  $k$  both children of node  $n$ , the entry of  $M$  in the row corresponding to  $k$  and the column corresponding to  $j$  is the aggregating of all endorsements from leaves in  $T_j$  to leaves in  $T_k$ . This entry is equal to the sum of all entries in the  $j$ th judgment corresponding to leaves of  $T_k$ .

To facilitate recursive implementation of our algorithm, we decided to localise the process of computing the local ranking for all the nodes while ignoring the links to and from the sub-tree. Localizing of marginal rank computations is motivated by the aforementioned "peer-review principle". We adopted QuickRank's definition of approximate endorsements by individuals outside the sub community. For instance, considering an individual  $i$  who is a member of community  $A$  and another individual  $j$  belonging to a different community  $B$ , An endorsement by  $i$  in  $A$  for  $j$  in  $B$  is as a component reflecting part of an endorsement by  $A$  of  $B$ . Thus, endorsements by  $A$  for individuals in  $B$  are aggregated into an endorsement by  $A$  of  $B$  by first scaling the endorsements from each  $i$  to each  $j$  by  $i$ 's local rank, and then adding the resulting weighted endorsements. Replacing  $j$  by any other  $j'$  in  $B$  does not change the aggregated endorsement. This is why the main endorsement is considered to be estimated. If the links starting at  $i$  are interpreted to mean  $i$ 's judgment, then this collection procedure is considered to be based on Bonacich's hypothesis. That is, obtaining endorsements of  $j \in B$  by  $A$  is followed by aggregating over all  $j \in B$  (to derive an endorsement of  $A$ ).

**Aggregation of Rankings and Links:** Based on QuickRank, the ranking of  $n$ 's  $m$  children into a single ranking corresponding to  $m$  can be achieved by calculating the means of the rankings  $r_1, \dots, r_m$  according to the weights indicated by the local ranking  $r$ . Concatenating the  $m$  rankings into a matrix  $Q = [r_1 \ \dots \ r_m]$ , the aggregation of ranking is simply expressed as  $Qr$ . Also corresponded with each child  $j$  of a collapsible node  $n$  are a judgment  $l_j$ , and a user bias  $b_j$ . These judgments and user bias are computed using the exact same techniques as the case of rankings.



Having described the major steps of our Enhanced Recursive Ranking Algorithm, we now present the algorithm in pseudo code (Table 3.1) (we refer the reader to Greenwald and Wicks (2006) for a detailed explanation on BaseRank procedure). In the pseudo code, weighted judgment is denoted as  $e_n$ , weighted bias as  $\beta_n$ . The matrices are denoted by  $P = [l_1 \ \cdots \ l_m]$  - judgment;  $Q = [r_1 \ \cdots \ r_m]$  - ranking;  $B = [\beta_1 \ \cdots \ \beta_m]$  - bias.

**Table 1. Enhanced Recursive Ranking Algorithm**

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*Algorithm: Enhanced Recursive Ranking (Collapsing node n)*

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- 1: **if** ( $n.isLeaf()$ ) **then**
- 2:     **return** ( $(n.getJudgment(), cn), Concat(\{n.getUserBias(), \beta_n\})$ );
- 3:     **Else**
- 4:          $m = n.numChildren()$
- 5:         **for** ( $j = 1$  to  $m$ ) **do**
- 6:              $(l_j, r_j, b_j) \leftarrow QuickRank(n.getChild(j))$
- 7:             **for** ( $k = 1$  to  $m$ ) **do**
- 8:                  $M_{kj} = Sum(l_j, n.getChild(k))$
- 9:             **end for**
- 10:         **end for**
- 11:          $P = [l_1 \ \cdots \ l_m]$
- 12:          $Q = [r_1 \ \cdots \ r_m]$
- 13:          $B = [\beta_1 \ \cdots \ \beta_m]$
- 14:          $r = BaseRank(M, n.getLocalPriorRanking())$
- 15:         **return** ( $(P_r, Q_r, B_r)$ );
- 16:     **end if**

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In the sample school structure described in previous section, using ERRA to compute the ranking of any department or faculty, we must compute the marginal (local) rankings of interior and posterior nodes (faculties and departments) and compute the cumulative local rankings using chain rule. Ranking techniques will be applied on each individual within the hierarchy. We begin by restricting the link graph a particular faculty, thereby constructing a local link subgraph including only the departments in that faculty. Then using BaseRank procedure, we compute the marginal ranking of the departments. Then we scale the links from this faculty's departments to the outside departments by the marginal ranks of the faculty's departments. Finally, the results are added to generate a resultant link which is represented from the current faculty to the outside connected departments. Furthermore, we progressed to recursively construct a marginal link sub graph, and generated corresponding ranking values for the faculties. The combination of the two local ranking for the departments in the faculty node produced one marginal ranking for all the departments in the faculty. Repeating this procedure for the other faculties, we obtain a single global ranking for each faculty. This process is repeated till the hierarchy is fully collapsed into a single a node.at this point, the ranking of all the departments in the school are generated.



Thus, obtained rank is the same as the QuickRank of the departments. We then scale the QuickRank using the pre-computed user bias aggregated at each node, to obtain the enhanced recursive ranking. If the user bias in all the nodes is not equal to 1, then the final ranking may differ from the QuickRank.

#### 4. NORMALIZATION OF RANKING CRITERIA

As stated in the description of the steps of the algorithm, it is required that the user bias be normalised. We explain why normalized bias value is needed, and the method of normalization. The third parameter in our algorithm, user bias, incorporates individualism into the ranking. The values of the user bias are computed from the user-defined search criteria. The order of preference of the search criteria (which may be specified explicitly by the user, or inferred implicitly from their profile) is important in the ranking of the results of the search. It is more convenient to allow users express these criteria in a form, most familiar to them. For instance, they may specify them as being ‘very important’, ‘important’, ‘not important’, etc. However, the quantification of these terms is arbitrary, and may not reflect the reality of the user’s intentions, as different users may understand these terms in different ways.

Another possible scheme is to allow them specify them on a scale of importance, for example, 0-10, 1-5, etc. Most surveys, with which people are accustomed, usually use such schemes. It makes sense, therefore, to adopt such an approach to specifying user-defined preferences. In order to make a direct comparison among criteria on their scale of relevance, it may be required that the cumulative sum of all the criteria equals a fixed number (possibly the maximum limit of the range). For instance, if we are using the 0-10 scale of relevance, it may be specified that the sum of the values, assigned by each user to the criteria must be equal to 10. Say, there are four criteria—newness, proximity, cheapness, ease of use. Then a user could assign any non-negative values to these criteria, as long as the sum equals 10. For example, one may assign 5 to newness, 3 to proximity, 0 to cheapness, 2 to ease of use; or 2.5 to all the criteria. We could then interpret the first example as the user considered newness to be of highest importance among the criteria., the next is proximity and then ease of use in that order, while the user does not consider cheapness important. In the second case, the user considers all the criteria to be equally important. Of course, while this requirement of constant cumulative sum enables us to quickly determine that order of preference of the criteria, as well as quantitatively determine the relative relevance of each, it is however, not very convenient for users to express their preferences that way. Users may know that some criteria are more important to them than others, but may not be able to quantitatively define how relatively important they are, that is, they are not able to say, whether criterion A is twice as important as criterion B. Moreover, such a scheme may not reflect the reality of user’s preferences.

As stated in the introduction to this chapter, a key question in this research work is determining a way to help users find what they intend, when they cannot specify in vivid terms exactly what they want. It is important, therefore, to allow users express their preferences in the form most familiar and convenient to them. To this end, we shall adopt one of the methods used for surveys—using a scale of preference to specify user search criteria. The scales are determined individually for each criterion, in such a way that, they make the most sense to the users. To illustrate, using our previous examples with four ranking criteria, we may adopt a scale of 0-10 for newness and cheapness, 1-5 for proximity, and 0-100 for ease of use. Of course, the same scale could be used for all the criteria, if it makes sense. However, the adoption of individual scales for the criteria, while very convenient, poses a problem - how can we determine the order of preferences, since different scales are used? To solve this problem, there is need to bring the values to the same scale for ease of comparison. This is achieved through normalization of the values.



There are various normalisation techniques in use today—min-max, cosine, non-monotonic, Z-score, etc. In order to select the best normalisation method to use, there is need to consider the purpose of such normalisation, the statistical properties of the data, and so on. Z-score normalization is useful if the distribution of the data is normal, which is not the case (or cannot be proven to be the case) for the values assigned the criteria. Local linear rescaling using quantile or percentile values alleviates the so-called ‘fat tail’ problem (Ready and Hu 1995), and therefore works well if there are outliers in the dataset, which is not the case here. For our purpose, we shall use the linear min-max normalization method. Since the linear min-max normalization does not distort the relative pairwise distances (because it is linear), it is the most appropriate for making our various criteria comparable without introducing additional errors. In min-max normalization, care must be taken to check for the  $min=max$  case, which luckily in our case, we do not need to worry about. It is computationally easy too. The formula for the min-max normalization is given by:

$$newX = \frac{\max - X}{\max - \min} \quad (3.1)$$

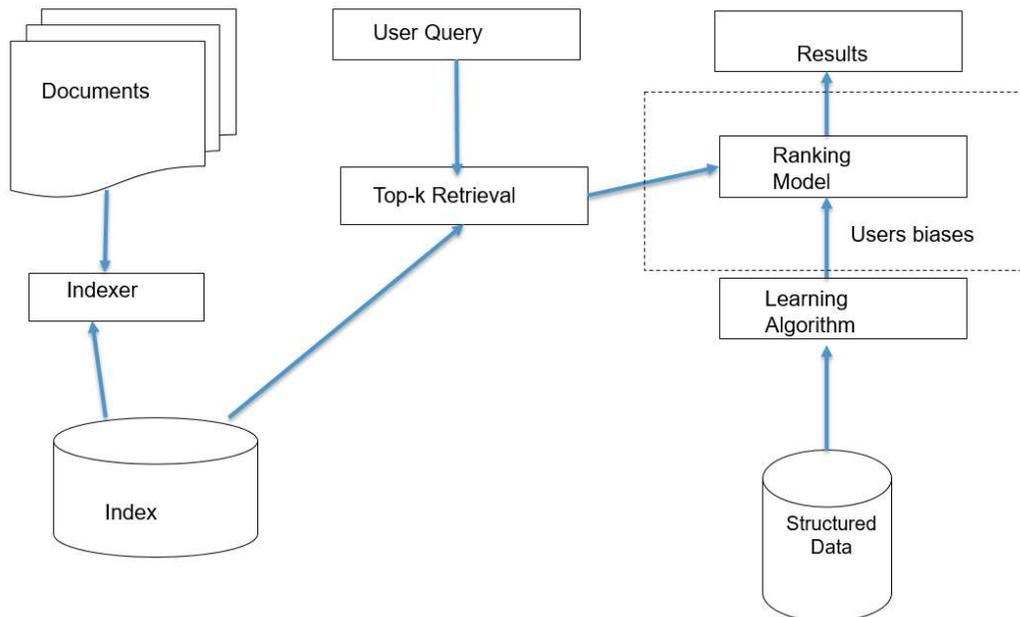
Where  $newX$  the normalized value of the criterion is,  $X$  is the user-specified value of the criterion,  $max$  refers to maximum limit of the scale,  $min$  represents minimum limit of scale. The min-max method normalizes each of the criteria in the range  $[0;1]$ . We shall interpret the value of 0 as meaning that the criterion is not important, while 1 signifies that the criterion is essential to the search. Larger values of the criteria indicate higher importance.

Our algorithm uses the min-max metric to normalize the ranking criteria. The purpose of normalizing the criteria is to be able to compare the order of preference of the ranking criteria, specified by the user. It also allows us to compare search results across similar data from different datasets. Using the min-max metric ensures that additional errors are not introduced by the normalisation step, as the relative distances between values are preserved. This method allows us to determine the importance of each of the ranking criteria in regards to the relationship between the QuickRank of the objects and their respective position in enhanced recursive ranking. The idea behind this normalisation metric is that most criteria are as important as the size of their normalised values, which in turn ensures that highly relevant results possess higher ranks in the generated list.

Finally, it is important to note that normalisation of the criteria is a pre-processing step, just like converting the link graph into a judgment graph. The user bias incorporated into our algorithm is actually not the user-specified values, but instead the normalized values. In this work the values of all the criteria are being normalized by using a min-max factor of the user-defined criteria, which removes the difficulty of the comparison of the relative importance of the ranking criteria, and allows us to compare search results across different datasets.

## 5. IMPLEMENTATION OF ENHANCED RECURSIVE RANK ALGORITHM

One of the deliverables of this research work is a software program that demonstrates our ranking algorithm. Here, we present explanation about the tools and techniques that were deployed for the implementation, special characteristics and limitations of the program. The structure of the software implementation of our enhanced ranking algorithm is shown on fig. 3.3 below. The components are discussed briefly.



**Fig. 2. Enhanced Recursive Ranking Model Implementation**

- i. **Indexer:** Our ranking algorithm enables users search for what they want, specifying the criteria and properties in order of preference. Of course, the traits will vary in importance to different users. A recursive search through a huge dataset places high demand on the processing power of the processing machine. It is therefore, required that the database be perfectly normalized and properly indexed. The task of indexing the records in the database is carried out by an indexer, inbuilt into the ranking system. Indexing makes the search easier, and ensures the fast operation of the database.
- ii. **User Query Handler:** To enable users input data, select search criteria, traits, qualities, etc., a robust and user-friendly interface is needed. The interface is developed in accordance with industrial standards. It receives user's query and link it to most relevant profile on the network. A **query processor** passes the queries through a web mining process, and returns the most relevant results.
- iii. **Database Manager:** A robust database is required to manage the huge volume of data. Owing to its ease of use, and free availability, MySQL database is used in the design of the ranking system.
- iv. **Web crawler:** A web crawler is used for the selection and gathering of the standardized dataset, which is used for the testing of the algorithm. From the web, it gathers relevant data, which forms the foundation of the database. The selected information will target people and their various unique characteristics, and will be stored in the database for referencing. The crawler is designed with the ability for data gathering from online social networks through their Application Programming



Interface (API). Using the APIs, it is possible to extract relevant information from the existing social network database, including data about the structure of the network. A popular example is Facebook API).

- v. **Ranker (or Ranking Model):** The ranker measures the importance of the search results returned, using the proposed Enhanced Recursive Ranking Algorithm. The ranker incorporates user bias in computing the ranks of the results.
- vi. **Retrieval Engine (Top-k Retrieval):** It performs lookups on the index tables against the search query.
- vii. **Learning Algorithm:** This implements a machine learning algorithm, which helps the system to adapt itself over time to rank results better, by analyzing previous data.
  - i. Our program is implemented using the Python and JavaScript computing language. The selection of these languages is based on the ease of programming, availability of computational libraries, as well as toolkits for GUI programming.

## 5. THE RANKING DISTRIBUTION

As stated previously, the ranking output by the algorithm is a probability distribution. From the statistical properties of this probability distribution, useful information about the dataset, and consequently the underlying hierarchical social network, can be extracted. Two of the most important properties are the **weighted mean** and **variance** of the distribution. **Weighted mean** enables us to calculate the estimated value of the rank given user bias. This is calculated by

$$\bar{x} = \frac{\sum_{i=1}^n w_i x_i}{\sum_{i=1}^n w_i} = \frac{w_1 x_1 + w_2 x_2 + \dots + w_n x_n}{w_1 + w_2 + \dots + w_n}$$

Where  $\{x_1, x_2, \dots, x_n\}$  is the dataset (ranks),  $\{w_1, w_2, \dots, w_n\}$  are the non-negative weights of the criteria. The weights are assigned according to the importance of the criteria at each node. It can be seen that criteria with a high weight contribute more to the weighted mean than do criteria with a low weight. We normalize the weights, such that their sum equals 1, that is  $\sum_{i=1}^n w_i = 1$ . Consequently, the formula above gets simplified to

$$\bar{x} = \sum_{i=1}^n w_i x_i = w_1 x_1 + w_2 x_2 + \dots + w_n x_n.$$

If equal normalised weights are assigned to all the criteria, then we get

$$\bar{x} = \frac{1}{n} \sum_{i=1}^n x_i$$

On the other hand, **variance** shows the ‘spread’ of the ranks. Higher variance signifies a wide difference in the relevance of the returned results, while lower variance means that the results are more or less on the same level of relevance. The variance of the distribution is computed by:



$$\text{Var}(X) = \frac{1}{n} \sum_{i=1}^n (x_i - \mu)^2$$

Where  $\mu = \frac{1}{n} \sum_{i=1}^n x_i$  is the expected value.

The weighted mean and variance are very significant, as they help us compare the degree of similarity between two standardized datasets (rankings) of the same network.

It is possible to determine the approximate **probability distribution of the rank**. In statistics there are numerous techniques (Ritzema 1994) for probability distribution fitting to a series of data, including parametric methods (Cramér 1946) and regression method. The parametric methods include; method of moments, method of L-moments (*Hosking 1990*), maximum likelihood method (*Aldrich 1997*). In this work, we use MATLAB's ALLFITDIST function, which fits all valid parametric probability distributions to data (MATLAB 2016).

Note that for non-parametric kernel-smoothing, FITDIST should be used directly instead. The probability distribution enables us to determine the dependence of the rank on different factors, from which useful information can be inferred.

The **correlation** between the original ranking (using QuickRank, or another ranking algorithm) and the individually biased ranking can be established. This correlation enables us to determine the influence of user bias in the ranking. This way, we could determine the efficiency of our algorithm in 'personalizing' search results based on user-defined preferences. Correlation can be computed using multiple methods. In this work, we adopt the Pearson Correlation Coefficient. The choice of method is based on the ease of computation, as well as the structure of the underlying test datasets. The Pearson Correlation Coefficient is given by:

$$\rho_{X,Y} = \frac{\text{cov}(X,Y)}{\sigma_X \sigma_Y} = \frac{E[(X - \mu_X)(Y - \mu_Y)]}{\sigma_X \sigma_Y} = \frac{E[XY] - E[X]E[Y]}{\sqrt{E[X^2] - (E[X])^2} \sqrt{E[Y^2] - (E[Y])^2}}$$

Where  $\text{cov}(X,Y)$  is the covariance,  $\sigma_X$  is the standard deviation of  $X$ ,  $\sigma_Y$  is the standard deviation of  $Y$ ,  $\mu_X$  is the mean of  $X$ ,  $\mu_Y$  is the mean of  $Y$ ,  $E$  is the expectation.

When applied to a sample, the Pearson Correlation Coefficient is given by:

$$r = r_{xy} = \frac{n \sum_{i=1}^n x_i y_i - \sum_{i=1}^n x_i \sum_{i=1}^n y_i}{\sqrt{n \sum_{i=1}^n x_i^2 - \left(\sum_{i=1}^n x_i\right)^2} \sqrt{n \sum_{i=1}^n y_i^2 - \left(\sum_{i=1}^n y_i\right)^2}} \quad (3.7)$$

for two datasets  $\{x_1, \dots, x_n\}$ , and  $\{y_1, \dots, y_n\}$ . To get the weighted correlation coefficient, we use the formula



$$\text{corr}(x, y; w) = \frac{\text{cov}(x, y; w)}{\sqrt{\text{cov}(x, x; w) \text{cov}(y, y; w)}} \quad (3.8)$$

where  $\text{cov}(x, y; w) = \frac{\sum_{i=1}^n w_i (x_i - m(x; w))(y_i - m(y; w))}{\sum_{i=1}^n w_i}$  is the weighted covariance;  $m(x; w) = \frac{\sum_{i=1}^n w_i x_i}{\sum_{i=1}^n w_i}$  is the

weighted mean.

In our case we apply formula (3.7) to the rankings obtained with ERRA, and the original rankings obtained through QuickRank or another algorithm. The value of  $r$  determines how strongly the introduction of user bias distorts the original ranking, how much ‘personalized’ by user criteria it becomes. Strong correlation coefficient can be interpreted to mean that user bias greatly influence the ranking result.

Another useful property is the **probability distribution of the search criteria**. It reflects how each criterion varies among users. The distribution shades light on the properties of the users, and could be used to train the system to predict and rank results better, using information inferred from the user’s profile. For this purpose, a learning algorithm unit is included in the program (see fig. 3.2). To determine the probability distribution of each of the criteria, we use the ALLFITDIST function from MATLAB as used above.

### 5.1. Generality and Limitations of the Enhanced Recursive Rank Algorithm

Having outlined in details the workings, implementation and the properties of the ERRA, it is necessary that we discuss the generality and limitations of the algorithm. As previously mentioned, the general principles of the proposed approach can be applied to any social network with a hierarchical structure. This implies that our algorithm has a wide area of applications, since a lot of search objects can be represented with a hierarchical structure. From this fact, however, springs a serious limitation – the algorithm cannot be used directly with networks, which are not hierarchical. However, it is possible to modify the algorithm to work on such non-hierarchical networks, or modify the structure of the network to suit the algorithm. We have, however, not yet investigated this possibility. Despite this limitation, the algorithm can be applied to existing search engines as a means of personalizing search results.

## 6. CONCLUSION

We have discussed about the proposition of a new algorithm for hierarchical social networks. The new algorithm named “Enhanced Recursive Ranking Algorithm” (ERRA) is a modification of the QuickRank algorithm to incorporate user bias. The theoretical framework for the implementation of the algorithm was discussed, as well as the statistical properties of the results. From the theoretical analysis, we demonstrated that the algorithm shows lots of promises in solving the research problems proposed. In the next chapter, experimental study of the algorithm using the software program developed will be carried out. This will help us determine the adequacy of the proposed solution.



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