# ON MOLECULAR GRAPHS WITH SMALLEST AND GREATEST ZEROTH-ORDER GENERAL RANDIĆ INDEX* 

Yumei Hu, Xueliang Li, Yongtang Shi, Tianyi Xu<br>Center for Combinatorics and LPMC<br>Nankai University, Tianjin 300071, P.R. China<br>Ivan Gutman<br>Faculty of Science, University of Kragujevac, Kragujevac<br>Serbia and Montenegro

(Received April 8, 2005)


#### Abstract

A molecular $(n, m)$-graph $G$ is a connected simple graph with $n$ vertices, $m$ edges and vertex degrees not exceeding 4. If $d(v)$ denotes the degree of the vertex $v$, then the zeroth-order general Randić index ${ }^{0} R_{\alpha}$ of the graph $G$ is defined as $\sum_{v \in V(G)} d(v)^{\alpha}$, where $\alpha$ is a pertinently chosen real number. We characterize, for any $\alpha$, the molecular $(n, m)$-graphs with smallest and greatest ${ }^{0} R_{\alpha}$.


[^0]
## 1 Introduction

The Randić (or connectivity) index was introduced by Randić in 1975 and is defined as [25]

$$
R=R(G)=\sum_{u v \in E(G)}(d(u) d(v))^{-1 / 2}
$$

where $d(u)$ denotes the degree of the vertex $u$ of the graph $G$, and $E(G)$ is the edge set of $G$. Randić himself demonstrated [25] that this index is well correlated with a variety of physico-chemical properties of various classes of organic compounds. Eventually, this structure-descriptor became one of the most popular topological indices to which two books $[15,17]$, several reviews $[7,24,26]$ and countless research papers are devoted.

Like other successful structure-descriptors, the Randić index received considerable attention also from mathematical chemists and mathematicians. In particular, bounds for $R$ and graphs extremal with regard to $R$ were extensively studied $[1,2,4,8,9,10$, $11,12]$.

In 1998 Bollobás and Erdös [3] generalized $R(G)$ by replacing the exponent $-1 / 2$ by an arbitrary real number $\alpha$. This graph invariant is called the general Randić index and will be denoted by $R_{\alpha}=R_{\alpha}(G)$. Li and Yang [20] studied $R_{\alpha}$ for general $n$-vertex graphs, obtained lower and upper bounds for it, and characterized the corresponding extremal graphs. Later $\mathrm{Hu}, \mathrm{Li}$ and Yuan $[13,14]$ determined the trees extremal with regard to $R_{\alpha}$, whereas Li, Wang and Wei [19] gave lower and upper bounds for $R_{\alpha}$ of molecular ( $n, m$ )-graphs. Other mathematical studies of the general Randić index are found in $[5,6]$.

The zeroth-order Randić index, conceived by Kier and Hall [16], is

$$
{ }^{0} R={ }^{0} R(G)=\sum_{v \in V(G)} d(v)^{-1 / 2}
$$

Eventually, Li and Zheng [22] defined the zeroth-order general Randić index of a graph $G$ as

$$
{ }^{0} R_{\alpha}={ }^{0} R_{\alpha}(G)=\sum_{v \in V(G)} d(v)^{\alpha}
$$

for any real number $\alpha$.
Pavlović determined the graphs with maximum ${ }^{0} R$-index [23]. Li et al. [18] investigated the same problem for the topological index $M_{1}(G)$, one of the Zagreb indices, that is defined as $M_{1}(G)=\sum_{v \in V(G)} d(v)^{2}$. (Evidently, $M_{1} \equiv{ }^{0} R_{\alpha}$ for $\alpha=+2$.) They obtained sufficient and necessary conditions under which ( $n, m$ )-graphs have minimum $M_{1}$, and a necessary condition for an $(n, m)$-graph having maximum $M_{1}(G) . \mathrm{Li}$ and Zhao [21] characterized trees with the first three smallest and largest zeroth-order general Randić index, with the exponent $\alpha$ being equal to $k,-k, 1 / k$, and $-1 / k$, where $k \geq 2$ is an integer.

In this paper we investigate the zeroth-order general Randić index of molecular ( $n, m$ )-graphs, i. e., connected simple graphs with $n$ vertices, $m$ edges and maximum vertex degree at most 4 . We characterize the molecular ( $n, m$ )-graphs with extremal (maximum or minimum) zeroth-order general Randić index.

First we need to introduce some notation.
Denote by $D(G)=\left[d_{1}, d_{2}, \cdots, d_{n}\right]$ the degree sequence of the graph $G$, where $d_{i}$ stands the degree of the $i$-th vertex of $G$, and $d_{1} \geq d_{2} \geq \cdots \geq d_{n}$.

If there is a graph $G$, such that $d_{i} \geq d_{j}+2$, let $G^{\prime}$ be the graph obtained from $G$ by replacing the pair $\left(d_{i}, d_{j}\right)$ by the pair $\left(d_{i}-1, d_{j}+1\right)$. In other words, if $D(G)=$ $\left[d_{1}, d_{2}, \cdots, d_{i-1}, d_{i}, d_{i+1}, \cdots, d_{j-1}, d_{j}, d_{j+1}, \cdots, d_{n}\right]$, then $D\left(G^{\prime}\right)=\left[d_{1}, d_{2}, \cdots, d_{i-1}, d_{i}-\right.$ $\left.1, d_{i+1}, \cdots, d_{j-1}, d_{j}+1, d_{j+1}, \cdots, d_{n}\right]$.

Note that if $\alpha=0$ then ${ }^{0} R_{\alpha}(G)=n$, and if $\alpha=1$ then ${ }^{0} R_{\alpha}(G)=2 m$. Therefore, in the following we always assume that $\alpha \neq 0,1$.

Lemma 1.1 For the two graphs $G$ and $G^{\prime}$, specified above, we have
(i) ${ }^{0} R_{\alpha}(G)>{ }^{0} R_{\alpha}\left(G^{\prime}\right)$ for $\alpha<0$ or $\alpha>1$
(ii) ${ }^{0} R_{\alpha}(G)<{ }^{0} R_{\alpha}\left(G^{\prime}\right)$ for $0<\alpha<1$.

Proof. Since ${ }^{0} R_{\alpha}(G)=\sum_{v \in V(G)} d(v)^{\alpha}$, we have

$$
\begin{aligned}
{ }^{0} R_{\alpha}(G)-{ }^{0} R_{\alpha}\left(G^{\prime}\right) & =d_{i}^{\alpha}+d_{j}^{\alpha}-\left(d_{i}-1\right)^{\alpha}-\left(d_{j}+1\right)^{\alpha} \\
& =\left[d_{i}^{\alpha}-\left(d_{i}-1\right)^{\alpha}\right]-\left[\left(d_{j}+1\right)^{\alpha}-d_{j}^{\alpha}\right] \\
& =\alpha\left(\xi_{1}^{\alpha-1}-\xi_{2}^{\alpha-1}\right)
\end{aligned}
$$

where $\xi_{1} \in\left(d_{i}-1, d_{i}\right)$, and $\xi_{2} \in\left(d_{j}, d_{j}+1\right)$. So, by $d_{i} \geq d_{j}+2$, we have $\xi_{1}>\xi_{2}$. Then ${ }^{0} R_{\alpha}(G)>{ }^{0} R_{\alpha}\left(G^{\prime}\right)$ for $\alpha<0$ or $\alpha>1$, whereas ${ }^{0} R_{\alpha}(G)<{ }^{0} R_{\alpha}\left(G^{\prime}\right)$ for $0<\alpha<1$.

## 2 Extremal molecular ( $n, m$ )-graphs

Denote by $n_{i}$ the number of vertices of degree $i$ in a molecular $(n, m)$-graph $G$. Then we have

$$
\begin{equation*}
{ }^{0} R_{\alpha}(G)=n_{1}+2^{\alpha} n_{2}+3^{\alpha} n_{3}+4^{\alpha} n_{4} \tag{2.1}
\end{equation*}
$$

Theorem 2.1 Let $C^{*}$ be a molecular ( $n, m$ )-graph with degree sequence $\left[d_{1}, d_{2}, \cdots, d_{n}\right]$, such that $\left|d_{i}-d_{j}\right| \leq 1$ for any $i \neq j$. Then for $\alpha<0$ or $\alpha>1, C^{*}$ has the minimum zeroth-order general Randić index among all molecular ( $n, m$ )-graphs, whereas for $0<$ $\alpha<1, C^{*}$ has the maximum zeroth-order general Randić index among all molecular ( $n, m$ )-graphs. Moreover,

$$
{ }^{0} R_{\alpha}\left(C^{*}\right)= \begin{cases}2+2^{\alpha}(n-2) & \text { if } m=n-1 \\ 2^{\alpha}(3 n-2 m)+3^{\alpha}(2 m-2 n) & \text { if } n \leq m \leq\lfloor 3 n / 2\rfloor \\ 3^{\alpha}(4 n-2 m)+4^{\alpha}(2 m-3 n) & \text { if }\lfloor 3 n / 2\rfloor<m \leq 2 n\end{cases}
$$

Proof. We only consider the case $0<\alpha<1$, because the proof for the other case is fully analogous. Let $G$ be a molecular graph and $D(G)=\left[d_{1}, d_{2}, \cdots, d_{n}\right]$. If $G \not \nexists C^{*}$, then there must exist a pair $\left(d_{i}, d_{j}\right)$ such that $d_{i} \geq d_{j}+2$. By Lemma 1.1, the graph $G^{\prime}$, obtained by replacing the pair $\left(d_{i}, d_{j}\right)$ by the pair $\left(d_{i}-1, d_{j}+1\right)$, has a greater ${ }^{0} R_{\alpha}$-value than $G$. Consequently, $G$ is not a molecular ( $n, m$ )-graph with maximum zeroth-order general Randić index.

To show the existence, we construct the extremal $(n, m)$-graph $C^{*}$ (with minimum ${ }^{0} R_{\alpha}$ for $\alpha<0$ or $\alpha>1$, and with maximum ${ }^{0} R_{\alpha}$ for $0<\alpha<1$ ) by adding edges one by one. First, we start from a tree. There must be at least two 1-degree vertices in a tree. By Lemma 1.1, there does not exist any 3-degree vertex, and so the extremal tree must be the path $P_{n}$. Next we add an edge joining the two leaves of the path. In this way the degrees of all vertices become equal to two, and then we get a cycle. We continue by adding edges one by one, so as to maximize the number of 3-degree vertices, until either there remain no 2-degree vertices, or remains exactly one. If more edges need to be added, then we first connect the 2-degree vertex (if such does exist) with a non-adjacent 3-degree vertex, and continue by connecting pairs of nonadjacent 3-degree vertices. The construction is shown in Figure 2.1.


Figure 2.1 Constructing molecular graphs with extremal zeroth-order general Randić index, according to Theorem 2.1.

Theorem 2.2 Let $G^{*}$ be a molecular ( $n, m$ )-graph with at most one vertex of degree 2 or 3. If one of the following conditions holds:
(I) $m=n-1$
(II) $m \geq n \geq 6$, for $n=6, m \geq 10$, and for $n=7, m \neq 8$
then for $\alpha<0$ or $\alpha>1, G^{*}$ has the maximum zeroth-order general Randić index among all molecular ( $n, m$ )-graphs, whereas for $0<\alpha<1$, the same graph has the minimum zeroth-order general Randić index among all molecular ( $n, m$ )-graphs. Moreover,

$$
{ }^{0} R_{\alpha}\left(G^{*}\right)= \begin{cases}(4 n-2 m) / 3+4^{\alpha}(2 m-n) / 3 & \text { if } 2 m-n \equiv 0(\bmod 3) \\ (4 n-2 m-2) / 3+2^{\alpha}+4^{\alpha}(2 m-n-1) / 3 & \text { if } 2 m-n \equiv 1(\bmod 3) \\ (4 n-2 m-1) / 3+3^{\alpha}+4^{\alpha}(2 m-n-2) / 3 & \text { if } 2 m-n \equiv 2(\bmod 3)\end{cases}
$$

Proof. Again, we only consider the case $0<\alpha<1$, because the proof for the other case is similar. Let $G^{\prime}$ be a molecular $(n, m)$-graph and $D\left(G^{\prime}\right)=\left[d_{1}, d_{2}, \cdots, d_{n}\right]$. Let $G^{\prime}$ possess two vertices of degree 2 or 3 , i. e., let there be a pair $\left(d_{i}, d_{j}\right)$, such that $3 \geq d_{i} \geq d_{j} \geq 2$. Then by Lemma 1.1, there is a graph $G$, obtained by replacing the pair $\left(d_{i}, d_{j}\right)$ by the pair $\left(d_{i}+1, d_{j}-1\right)$, that has a smaller ${ }^{0} R_{\alpha}$-value than $G^{\prime}$. Repeating the above operation until there is no pair $\left(d_{i}, d_{j}\right)$, such that $3 \geq d_{i} \geq d_{j} \geq 2$, we arrive at $G^{*}$ with minimum zeroth-order general Randić index. In view of (2.1), for $G^{*}$ we have

$$
\begin{cases}n_{1}+n_{2}+n_{3}+n_{4} & =n \\ n_{1}+2 n_{2}+3 n_{3}+4 n_{4} & =2 m \\ n_{2}+n_{3} & \leq 1\end{cases}
$$

From the above equations, we have one of the following three options:
(1) $n_{2}=n_{3}=0$, implying $n_{1}=(4 n-2 m) / 3, n_{4}=(2 m-n) / 3$, and $2 m-n \equiv$ $0(\bmod 3)$
(2) $n_{2}=1, n_{3}=0$, implying $n_{1}=(4 n-2 m-2) / 3, n_{4}=(2 m-n-1) / 3$, and $2 m-n \equiv 1(\bmod 3)$
(3) $n_{2}=0, n_{3}=1$, implying $n_{1}=(4 n-2 m-1) / 3, n_{4}=(2 m-n-2) / 3$, and $2 m-n \equiv 2(\bmod 3)$.

In order to show the existence, we construct $G^{*}$ by distinguishing the following cases:
(I) $m=n-1$, i. e., $G^{*}$ is a tree.
(I.1) If $2 m-n \equiv 0(\bmod 3)$, we first construct a path with $n_{4}$ vertices, and then add $n_{1}$ pendent vertices, taking care that no vertex gets degree greater than 4.
(I.2) If $2 m-n \equiv 1(\bmod 3)$, we first construct a path with $n_{4}$ vertices, then add $n_{1}$ pendent vertices, taking care that no vertex gets degree greater than 4 , and finally subdivide an edge by inserting to it a vertex of degree 2 .
(I.3) If $2 m-n \equiv 2(\bmod 3)$, we first construct a path with $n_{4}+1$ vertices, and then add $n_{1}$ pendent vertices, taking care that no vertex gets degree greater than 4.
(II) $m \geq n \geq 6$, for $n=6$ and $m \geq 10$, or for $n=7$ and $m \neq 8$.

In Figure 2.2 we show one of the possible graphs $G^{*}$ for $n_{4} \leq 4$, that is for
$\lfloor(2 m-n) / 3\rfloor \leq 4$. For $n_{4} \geq 5$, we construct $G^{*}$ as follows:
(II.1) If $2 m-n \equiv 0(\bmod 3)$, we first construct a 4 -regular graph on $n_{4}$ vertices, then delete $n_{1} / 2$ edges from it, and then add $n_{1}$ pendent vertices, taking care that no vertex gets degree greater than 4 .
(II.2) If $2 m-n \equiv 1(\bmod 3)$, we first construct a 4 -regular graph on $n_{4}$ vertices, then delete $n_{1} / 2$ edges from it, then add $n_{1}$ pendent vertices, taking care that no vertex gets degree greater than 4 , and finally subdivide an edge inserting to it a vertex of degree 2 .
(II.3) If $2 m-n \equiv 2(\bmod 3)$, we first construct a 4 -regular graph on $n_{4}+1$ vertices, then delete $\left(n_{1}+1\right) / 2$ edges from it, and then add $n_{1}$ pendent vertices, taking care that no vertex gets degree greater than 4 .
This completes the proof.
Note that Theorem 2.2 holds under the conditions $m=n-1$, or $n=6$ and $m \geq 10$, or $n=7$ and $m \neq 8$, or $m \geq n \geq 8$, since for the other pairs of $n$ and $m$ the extremal degree sequences obtained in Theorem 2.2 are not graphic. It is easy to check that for $n=1,2,3$, and for $n \geq 4$ and $m=\binom{n}{2}-1$ or $m=\binom{n}{2}$ the ( $n, m$ )-graph is unique. For $n=4$ and $m=4$, or $n=5$ and $5 \leq m \leq 8$, or $n=6$ and $6 \leq m \leq 9$, or $n=7$ and $m=8$ we can characterize the extremal graphs by examining all possible degree sequences. These extremal graphs are depicted in Figure 2.3 (minimum ones for $0<\alpha<1$, maximum ones for $\alpha<0$ or $\alpha>1$, except for $n=5$ and $m=7$, in which case $(a)$ is the minimum graph for $0<\alpha<1$ and maximum graph for $\alpha<0$ or $1<\alpha<2$, and (b) is the maximum graph for $\alpha \geq 2$ ).

$n=6, \quad m=10$

$\mathrm{n}=7, \quad \mathrm{~m}=7$

$n=7, \quad m=9$

$\mathrm{n}=7, \quad \mathrm{~m}=10$

$\mathrm{n}=8, \quad \mathrm{~m}=8$

$\mathrm{n}=8, \quad \mathrm{~m}=9$

$n=8, \quad m=10$

$n=8, \quad m=11$

$n=9, \quad m=9$

$n=10, \quad m=12$

$\mathrm{n}=11, \quad \mathrm{~m}=11$

$\mathrm{n}=11, \mathrm{~m}=12$



$\mathrm{n}=13, \quad \mathrm{~m}=13$

$\mathrm{n}=14, \quad \mathrm{~m}=14$

Figure 2.2 Molecular graphs with extremal zeroth-order general Randić index, having four or fewer vertices of degree 4.


Figure 2.3 Some graphs with extremal zeroth-order general Randić index; for details see text.

## References

[1] O. Araujo, J. A. de la Peña, Some bounds for the connectivity index of a chemical graph, J. Chem. Inf. Comput. Sci. 38 (1998) 827-831.
[2] O. Araujo, J. Rada, Randić index and lexicographic order, J. Math. Chem. 27 (2000) 201-212.
[3] B. Bollobás, P. Erdös, Graphs of extremal weights, Ars Combin. 50 (1998) 225233.
[4] G. Caporossi, I. Gutman, P. Hansen, L. Pavlović, Graphs with maximum connectivity index, Comput. Biol. Chem. 27 (2003) 85-90.
[5] L. H. Clark, J.W. Moon, On the general Randić index for certain families of trees, Ars Combin. 54 (2000) 223-235.
[6] M. Fischermann, A. Hoffmann, D. Rautenbach, L. Volkmann, A linear-programming approach to the generalized Randić index, Discr. Appl. Math. 128 (2003) 375-385.
[7] S. P. Gupta, Quantitative structure-activity relationship studies on local anesthetics, Chem. Rev. 91 (1991) 1109-1119.
[8] I. Gutman, O. Araujo, D. A. Morales, Bounds for the Randić connectivity index, J. Chem. Inf. Comput. Sci. 40 (2000) 572-579.
[9] I. Gutman, O. Araujo, D. A. Morales, Estimating the connectivity index of a saturated hydrocarbon, Indian J. Chem. 39A (2000) 381-385.
[10] I. Gutman, O. Miljković, Molecules with smallest connectivity indices, MATCH Commun. Math. Comput. Chem. 41 (2000) 57-70.
[11] I. Gutman, O. Miljković, G. Caporossi, P. Hansen, Alkanes with small and large Randić connectivity index Chem. Phys. Lett. 306 (1999) 366-372.
[12] P. Hansen, H. Mélot, Variable neighborhood search for extremal graphs 6: Analyzing bounds for the connectivity index, J. Chem. Inf. Comput. Sci. 43 (2003) 1-14.
[13] Y. Hu, X. Li, Y. Yuan, Trees with minimum general Randić index, MATCH Commun. Math. Comput. Chem. 52 (2004) 119-128.
[14] Y. Hu, X. Li, Y. Yuan, Trees with maximum general Randić index, MATCH Commun. Math. Comput. Chem. 52 (2004) 129-146.
[15] L. B. Kier, L. H. Hall, Molecular Connectivity in Chemistry and Drug Research, Academic Press, New York, 1976.
[16] L. B. Kier, L. H. Hall, The nature of structure-activity relationships and their relation to molecular connectivity, Europ. J. Med. Chem. 12 (1977) 307-312.
[17] L. B. Kier, L. H. Hall, Molecular Connectivity in Structure-Analysis, Research Studies Press, Wiley, Chichester, UK, 1986.
[18] R. Lang, X. Li, S. Zhang, Inverse problem for Zagreb index of molecular graphs, Appl. Math. J. Chinese Univ. A18 (2003) 487-493 (in Chinese).
[19] X. Li, X. Q. Wang, B. Wei, On the lower and upper bounds for general Randic index of chemical ( $n, m$ )-graphs, MATCH Commun. Math. Comput. Chem. 52 (2004) 157-166.
[20] X. Li, Y. Yang, Sharp bounds for the general Randic index, MATCH Commun. Math. Comput. Chem. 51 (2004) 155-166.
[21] X. Li, H. Zhao, Trees with the first three smallest and largest generalized toplogical indeices, MATCH Commun. Math. Comput. Chem. 50 (2004) 57-62.
[22] X. Li, J. Zheng, A unified approach to the extremal trees for different indices, MATCH Commun. Math. Comput. Chem. 54 (2005) 195-208.
[23] L. Pavlović, Maximal value of the zeroth-order Randić index, Discr. Appl. Math. 127 (2003) 615-626.
[24] L. Pogliani, From molecular connectivity indices to semiempirical connectivity terms: Recent trends in graph theoretical descriptors, Chem. Rev. 100 (2000) 3827-3858.
[25] M. Randić, On characterization of molecular branching, J. Am. Chem. Soc. 97 (1975) 6609-6615.
[26] M. Randić, The connectivity index 25 years after, J. Mol. Graphics Modell. 20 (2001) 19-35.


[^0]:    *Supported by the National Science Foundation of China.

