

On the SD- polynomial and SD- index of an infinite class of “Armchair Polyhex Nanotubes”

Mohammad Reza Farahani

Department of Applied Mathematics, Iran University of Science and Technology (IUST),
Narmak, Tehran 16844, Iran

E-mail address: Mr_Farahani@Mathdep.iust.ac.ir , mrfarahani88@gmail.com

ABSTRACT

Let G be a simple connected graph with the vertex set $V = V(G)$ and the edge set $E = E(G)$, without loops and multiple edges. For counting qoc strips in G , Diudea introduced the Ω -polynomial of G and was defined as $\Omega(G, x) = \sum_{i=1}^k x^{c_i}$, where C_1, C_2, \dots, C_k be the “opposite edge strips” ops of G and $c_i = |C_i|$ ($i = 1, 2, \dots, k$). One can obtain the Sd -polynomial by replacing x^c with $x^{|E(G)|-c}$ in Ω -polynomial. Then the Sd -index will be the first derivative of $Sd(x)$ evaluated at $x = 1$. In this paper we compute the Sd -polynomial and Sd -index of an infinite class of “Armchair Polyhex Nanotubes”.

Keywords: Omega and Sadhana polynomial; Sadhana index; Armchair Polyhex Nanotubes and Nanotori

1. INTRODUCTION

By a graph G means a pair $G = (V, E)$ in which $V = V(G)$ and $E = E(G)$ denote to the set of vertices and edges, respectively. A chemical graph is a graph theoretical representation of a molecule whose vertices correspond to the atoms of the compound and edges correspond to chemical bonds. For two vertices x and y belong to V , x is adjacent to y if and only if $xy \in E(G)$. In a connected graph, there is a path between every pair (x, y) of its vertices. The distance $d(x, y)$ between vertices/atoms x and y ($x, y \in V(G)$) is defined as the length of a shortest path between x and y . Two edges $e = uv$ and $f = xy$ of G are called co-distant, “ e co f ”, if and only if they obey the following relation for a non-negative integer d : [1]

$$d(v, x) = d(v, y) + 1 = d(u, x) + 1 = d(u, y) = d$$

For some edges of G there are the following relations satisfied [1,2]:

$$e \text{ co } e$$

$$e \text{ co } f \Leftrightarrow f \text{ co } e$$

$$e \text{ co } f \ \& \ f \text{ co } h \Rightarrow e \text{ co } h$$

though the last relation is not always valid. In other words, the relation “co” is reflexive and symmetric but it is not necessary to be transitive. Set $C(e) := \{f \in E(G), \mid e \text{ co } f\}$, denote the subset of edges in G , co-distant to the edge e . If the relation “co” is transitive on $C(e)$ then $C(e)$ is called an *orthogonal cut* (denoted by *oc*) of G . The graph G is called *co-graph* if and only if the edge set $E(G)$ a union of disjoint orthogonal cuts: $E(G) = \bigcup_{i=1}^k C_i$ and $C_i \cap C_j = \emptyset$, for $i \neq j$ and $i, j = 1, 2, \dots, k$.

If any two consecutive edges of an edge-cut sequence are topologically parallel within the same face of the covering, such a sequence is called a *quasi-orthogonal cut qoc* strip. For counting “opposite edge strips” qocs C_i of $E(G)$ ($i, j = 1, 2, \dots, k$), *M.V. Diudea* introduced the Ω -polynomial of G [3-11] and was defined as $\Omega(G, x) = \sum_{i=1}^k x^{c_i}$, where c_i 's is the size of opposite edge strips ($= |C_i|$ ($i = 1, 2, \dots, k$)).

It is easy to see that the first derivative of Omega polynomial $\Omega(G, x)$ (in $x = 1$) equals the number of edges in the graph

$$\Omega'(G, x) = \sum_{i=1}^k c_i = \sum_{i=1}^k |C_i| = |E(G)|$$

Another polynomial also related to the *ops* in G was introduced by *Ashrafi* and co-authors [12] in 2008, that counting the non-opposite edges is the *Sadhana* polynomial $Sd(G, x)$ defined as:

$$Sd(G, x) = \sum_{i=1}^k x^{|E(G)|-c_i}$$

The *Sadhana* index $Sd(G)$ for counting *qoc* strips in G was defined by *Khadikar et. al* [13,14] as first derivative of *sadhana* polynomial evaluated at $x = 1$ [13-18]

$$Sd(G) = Sd'(G, x) = \sum_{i=1}^k (|E(G)| - c_i)$$

By definition of Ω -polynomial, one can obtain the Sd -polynomial by replacing x^c with $x^{|E(G)|-c}$ in Ω -polynomial.

In chemical, physics and nano sciences, we have the appealing structure, especially symmetric structure with chemical constitution purporting. Carbon exists in several forms in nature. One is the so-called nanotube which was discovered for the first time in 1991 [19,20]. One of the nanotube is *Polyhex Nanotubes*, that the structure of polyhex nanotubes is consisting of the cycles with length six C_6 in columns.

Since polyhex nanotubes have more practical in the chemical, physics and nano science, in this paper we focus on its structure and by using definition of Sd -polynomial and Sd -index, we compute these topological polynomial and index for an infinite class of Nano-structure “*Armchair Polyhex Nanotubes TUAC₆*”, depicted in Figure 1.

Throughout this paper our notation is standard and mainly taken from standard book of graph theory such as [21-25].

2. RESULTS AND DISCUSSION

In this section we compute the Sd -polynomial and Sd -index of a family of *Polyhex Nanotubes*. In Figure 1, one can see that the 3-dimensional and 2-dimensional graph of Armchair polyhex nanotubes $TUAC_6[m,n]$, where m,n are the numbers of rows/columns of hexagon (C_6) in 2-dimensional perception $TUAC_6[m,n]$. In a series of papers [26-36], some properties and applications and more historical details of nanotubes are presented and studied.

By these terminologies and from Figure 1, we will have the following results for *Armchair Polyhex Nanotubes* $TUAC_6$.

Theorem 1.

$\forall m,n \in N$ let $G = TUAC_6[m,n]$ be the Armchair polyhex nanotubes, then the Sd -polynomial and Sd -index of G are equal to

$$Sd(TUAC_6[m,n], x) = 2mx^{6mn+4m-n-1} + 2mx^{6mn+4m-2n-1}$$

and

$$Sd(TUAC_6[m,n]) = 24m^2n + 16m^2 - 6mn - 4m$$

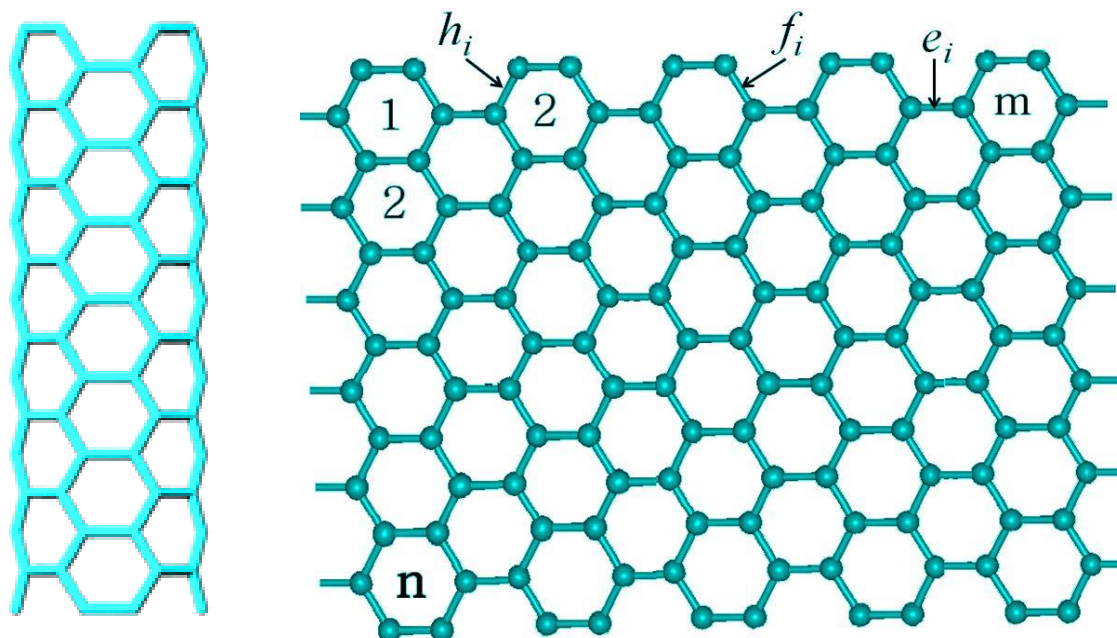


Fig. 1. A 3-dimensional (left) and 2-dimensional (right) lattices of Armchair Polyhex Nanotubes $TUAC_6[m,n]$.

Proof.

Consider the Armchair polyhex nanotubes $G = TUAC_6[m, n]$ ($m, n \in \mathbb{N}$) (Figure 1). Let m, n , $|V(G)|$ and $|E(G)|$ be the hexagons in rows/columns, number of vertices/carbon atoms and edges/chemical bonds of G . Then one can see that $|V(G)| = 4m(n + 1)$ and $|E(G)| = 6mn + 4m$. Now, if we denote all horizontal edge in i^{th} column by e_i and all left (or right) oblique edges in i^{th} column by f_i (or h_i), then it is easy to see that for all *quasi-orthogonal cuts* C_1, C_2, \dots, C_{2m} , $C_i = C(e_i)$ and also for all *quasi-orthogonal cuts* $C_{2m+1}, C_{2m+2}, \dots, C_{3m}$, $C_{2m+j} = C(f_j)$ and alternatively, for all *qocs* $C_{3m+1}, C_{3m+2}, \dots, C_{4m}$, $C_{3m+l} = C(h_l)$.

Now by according to Figure 1, one can see that $\forall i=1, 2, \dots, 2m: c_i = n + 1$ and $\forall j = 1, 2, \dots, m: c_{2m+j} = c_{3m+j} = 2n + 1$.

Thus, *Sd*-polynomial of Armchair polyhex nanotubes $G = TUAC_6[m, n]$ is equal to

$$\begin{aligned} Sd(TUAC_6[m, n], x) &= \sum_{i=1}^k x^{|E(G)|-c_i} \\ &= 2m \times x^{6mn+4m-n-1} + m \times x^{6mn+4m-2n-1} + m \times x^{6mn+4m-2n-1} \end{aligned}$$

The *Sd*-polynomial of G implies that the *Sd*-index of $TUAC_6[m, n]$ is equal to

$$\begin{aligned} Sd(TUAC_6[m, n]) &= Sd'(TUAC_6[m, n], x) = \left. \frac{\partial Sd(TUAC_6[m, n], x)}{\partial x} \right|_{x=1} \\ &= 2m \times (6mn + 4m - n - 1) + m \times (6mn + 4m - 2n - 1) + m \times (6mn \\ &\quad + 4m - 2n - 1) = 24m^2n + 16m^2 - 6mn - 4m \end{aligned}$$

Here, the proof is completed.

3. CONCLUSION

In this paper, we obtained the Sadhana polynomial and Sadhana index of Armchair Polyhex Nanotubes and Nanotori for the first time.

References

- [1] P.E. John, A.E. Vizitui, S. Cigher, and M.V. Diudea, *MATCH Commun. Math. Comput. Chem.* 57 (2007) 479-484.
- [2] M.V. Diudea, S. Cigher and P.E. John. Omega and Related Counting Polynomials. *MATCH Commun. Math. Comput. Chem.* 60 (2008) 237-250.
- [3] M.V. Diudea, S. Cigher, and P.E. John, *MATCH Commun. Math. Comput. Chem.* 60 (2008) 237-250.
- [4] M.V. Diudea, I. Gutman, L. Jäntschi, *Molecular Topology*, NOVA, New York, 2002.
- [5] M.V. Diudea, *Carpath. J. Math.* 22 (2006) 43-47.

-
- [6] M.V. Diudea, S. Cigher, A.E. Vizitiu, O. Ursu, P.E. John, *Croat. Chem. Acta* 79(3) (2006) 445-448.
- [7] P.E. John, A.E. Vizitiu, S. Cigher, M.V. Diudea, *MATCH Commun. Math. Comput. Chem.* 57 (2007) 479.
- [8] A.E. Vizitiu, S. Cigher, M.V. Diudea, M. S. Florescu, *MATCH Commun. Math. Comput. Chem.* 57(2) (2007) 479-484.
- [9] A.R. Ashrafi, M. Jalali, M. Ghorbani, M.V. Diudea, *MATCH, Commun. Math. Comput. Chem.* 60 (2008) 905-916.
- [10] M.V. Diudea, A. Ilić, *Carpath. J. Math.* 20(1) (2009) 177-185.
- [11] M.V. Diudea, *MATCH Commun. Math. Comput. Chem.* 64 (2010) 569.
- [12] A.R. Ashrafi, M. Ghorbani, M. Jalali, *Int. J. Chem.* 47A(4) (2008) 535-537.
- [13] P.V. Khadikar, S. Joshi, A.V. Bajaj, D. Mandloi, *Med. Chem. Lett.* 14 (2004) 1187-1191.
- [14] P.V. Khadikar, V.K. Agrawal, S. Karmarkar, *Bioorg. Med. Chem.* 2(10) (2002) 3499-3507.
- [15] P.V. Khadikar, D. Mandoli, *Bioinformatics Trends* 1 (2006) 51-63.
- [16] Khadikar, S. Singh, M. Jaiswal, D. Mandoli, *Bioorg. Med. Chem. Lett.* 14 (2004) 4795-4801.
- [17] P.V. Khadikar, J. Singh, M. Ingle, *J. Math. Chem.* 42 (2007) 433-446.
- [18] H. Mesgarani, M. Ghorbani, *Adv. Mater. Rapid Commun.* 4(11) (2010) 1863-1865.
- [19] S. Iijima, *Nature* 354 (1991) 56.
- [20] D. S. Bethune, C. H. Kiang, M. S. Devries, G. Gorman, R. Savoy, J. Vazquez, A. Beyers, *IBID* 363 (1993) 605.
- [21] F. Harary, *Graph Theory*, Addison-Wesley, Reading, MA, (1969).
- [22] B. West, *Introduction to graph theory*. Prentice Hall of India, (2003).
- [23] N. Trinajstić, *Chemical Graph Theory*, (second ed.) CRC Press, Boca Raton, FL, (1992).
- [24] R. Todeschini, V. Consonni, *Handbook of Molecular Descriptors*, Weinheim, Wiley-VCH, (2000).
- [25] N. Trinajstić, I. Gutman, *Croat. Chem. Acta* 75 (2002) 329-356.
- [26] I. Gutman, S. Klavžar, *ACH Models Chem* 133 (1996) 389-399.
- [27] M.V. Diudea, *MATCH, Commun. Math. Comput. Chem.* 45 (2002) 109-122.
- [28] A.R. Ashrafi, G. R. Vakili-Nezhaad, *Journal of Physics: Conference Series* 29 (2006) 181-184.
- [29] S. Yousefi, H. Yousefi-Azari, A.R. Ashrafi, M. H. Khalifeh, *JUST* 33(3) (2008) 7-11.
- [30] A. Iranmanesh, Y. Alizadeh, *Digest. J. Nanomater. Bios* 4 (2009) 607-611.
- [31] H. Shabani, A.R. Ashrafi, *Digest. J. Nanomater. Bios* 4 (2009) 423-428.

-
- [32] S. Alikhani, M.A. Iranmanesh, *Digest. J. Nanomater. Bios* 5 (2010) 1-7.
- [33] M.R. Farahani, *Romanian Academy Series B Chemistry* 15(1) (2013) 3-6.
- [34] M.R. Farahani, *Journal of Advances in Physics* 3(1) (2013) 191-196.
- [35] M.R. Farahani, *Acta Chim. Slov.* 59 (2012) 779-783.
- [36] M.R. Farahani. Computing GA_5 Index of Armchair Polyhex Nanotube, *Le Matematiche*. In press (2014).
- [37] Mohammad Reza Farahani, *International Letters of Chemistry, Physics and Astronomy* 11(1) (2014) 74-80.

(Received 20 March 2014; accepted 26 March 2014)