

# A highly correlated topological index for polyacenes

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## Abstract

The most used molecular graph descriptors in establishing Quantitative structure-property relationships (QSPRS) and Quantitative structure-activity relationships (QSARS) are topological indices. Molecular descriptors are normally chosen based on their ability to give good results in statistical models. In this paper we introduce a set of five new indices (Kekule indices)  $K$ ,  $K_1$ ,  $K_2$ ,  $K_3$ ,  $K_4$  and we establish that the Kekule index ( $K$ ) has excellent correlation ( $r = 0.99999997250969$ ) with  $\log p$  values in case of polyacenes.

**Keywords:** Kekule index, Polyacenes.

## INTRODUCTION

A topological index of a chemical compound is an integer, derived following a certain rule, which can be used to characterize the chemical compound. The first topological index is Wiener index introduced by Harold Wiener in 1947 to demonstrate its relation with physicochemical properties of alkanes, alcohols and amines. Ever since it is known that topological indices can be used to establish Quantitative structure - property relationships (QSPRS) and Quantitative structure - activity relationships (QSARS) in pharmacology, researchers are pursuing several investigations to find topological indices having correlation one or nearer to one with physicochemical properties of organic compounds. In this paper we introduce a new index  $K$  (Kekule index) having correlation ( $r = 0.99999997250969$ ) with  $\log p$  values of polyacenes.

## DEFINITIONS

In this section we define five new topological indices and explain the procedure of computation.

**Kekule index:** The Kekule index of a graph  $G = (V, E)$  is defined as  $K(G) = \sum_{e=uv \in E(G)} W(e)$  where  $W(e) = |i - j|$ ,  $i, j$  are the degrees of the vertices  $u$  and  $v$  in  $G$ .

**$K_1$  index:** The  $K_1$  index of a graph  $G = (V, E)$  is defined as  $K_1(G) = \sqrt{\sum_{e=uv \in E(G)} W(e)}$  where  $W(e) = |i - j|^2$ ,  $i, j$  are the degrees of the vertices  $u$  and  $v$  in  $G$ .

**$K_2$  index:** The  $K_2$  index of a graph  $G = (V, E)$  is defined as  $K_2(G)$

$= \sum_{e=uv \in E(G)} W(e)$  where  $W(v) =$  sum of the degrees of the neighboring vertices in  $G$ .

**$K_3$  index:** The  $K_3$  index of a graph  $G = (V, E)$  is defined as  $K_3(G) = \sum_{e=uv \in E(G)} W(e)$  where  $W(v) =$  product of the degrees of the neighboring vertices in  $G$ .

**$K_4$  index:** The  $K_4$  index of a graph  $G = (V, E)$  is defined as  $K_4(G) = \sum_{e=uv \in E(G)} W(e)$  where  $W(v) = \max\{d(v, v_i) / v_i \in V(G)\}$ . Where  $d(v, v_i)$  is the maximum distance between  $v$  and  $v_i$ .

We compute these indices considering the chemical graph of the compound Naphthalene given below.

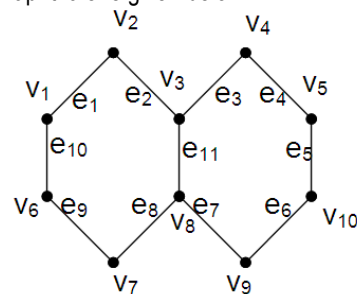


Fig 1. Graph of Naphthalene

Calculation of  $K$  index:

$W(e_1) = |2-2| = 0$ ,  $W(e_2) = |2-3| = 1$ ,  $W(e_3) = |3-2| = 1$ ,  $W(e_4) = |2-2| = 0$ ,  
 $W(e_5) = |2-2| = 0$ ,  $W(e_6) = |2-2| = 0$ ,  
 $W(e_7) = |3-2| = 1$ ,  $W(e_8) = |3-2| = 1$ ,  $W(e_9) = |2-2| = 0$ ,  $W(e_{10}) = |2-2| = 0$ ,  
 $W(e_{11}) = |3-3| = 0$

Therefore  $K(G) = 1 + 1 + 1 + 1 = 4$ .

Calculation of  $K_1$  index:

$W(e_1) = |2-2|^2 = 0$ ,  $W(e_2) = |2-3|^2 = 1$ ,  $W(e_3) = |3-2|^2 = 1$ ,  $W(e_4) = |2-2|^2 = 0$ ,  
 $W(e_5) = |2-2|^2 = 0$ ,  
 $W(e_6) = |2-2|^2 = 0$ ,  $W(e_7) = |3-2|^2 = 1$ ,  $W(e_8) = |3-2|^2 = 1$ ,  $W(e_9) = |2-2|^2 = 0$ ,  
 $W(e_{10}) = |2-2|^2 = 0$ ,  
 $W(e_{11}) = |3-3|^2 = 0$

Therefore  $K_1(G) = \sqrt{1+1+1+1} = 2$ .

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Calculation of  $K_2$  index:

$$W(v_1)= 2+2=4, W(v_2)= 2+3=5, W(v_3)= 2+3+2=7, W(v_4)= 3+2=5, \\ W(v_5)= 2+2=4, \\ W(v_6)= 2+2=4, W(v_7)= 2+3=5, W(v_8)= 2+2+3=7, W(v_9)= 3+2=5, \\ W(v_{10})= 2+2=4, \\ \text{Therefore } K_2(G) = 4+5+7+5+4+4+5+7+5+4=50.$$

Calculation of  $K_3$  index:

$$W(v_1)= 2 \times 2=4, W(v_2)= 2 \times 3=6, W(v_3)= 2 \times 3 \times 2=12, W(v_4)= \\ 3 \times 2=6, W(v_5)= 2 \times 2=4, \\ W(v_6)= 2 \times 2=4, W(v_7)= 2 \times 3=6, W(v_8)= 2 \times 2 \times 3=12, W(v_9)= \\ 3 \times 2=6, W(v_{10})= 2 \times 2=4, \\ \text{Therefore } K_3(G) = 4+6+12+6+4+4+6+12+6+4=64.$$

Calculation of  $K_4$  index:

$$W(v_1) = 5, W(v_2) = 4, W(v_3) = 3, W(v_4) = 4, W(v_5) = 5, W(v_6) = 5, \\ W(v_7) = 4, \\ W(v_8) = 3, W(v_9) = 4, W(v_{10}) = 5, \\ \text{Therefore } K_4(G)=5+4+3+4+5+5+4+3+4+5=42.$$

**NOTATION**

The molecular graph of polyacenes is a chain of hexagons arranged linearly.

For convenience we adopt the following notation:

Let  $L(a)$  be the graph consisting of  $a$  hexagons in **one** row as shown in the figure below. Here  $a$  is a positive integer.

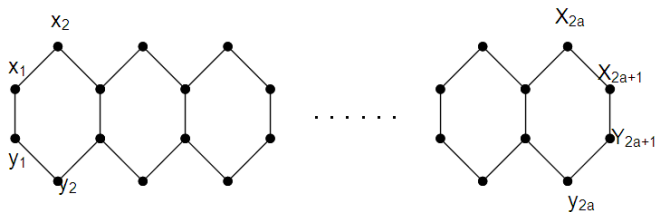


Fig 2. Graph of polyacene with a hexagons in 1 row

**MAIN RESULTS**

Throughout the paper we take

$$A = \{x_1, x_2, \dots, x_{2a+1}\}, B = \{y_1, y_2, \dots, y_{2a+1}\}, \\ E_1 = \{x_i, x_{i+1} / i = 1, 2, \dots, 2a\}, E_2 = \{y_i, y_{i+1} / i = 1, 2, \dots, 2a\}, E_3 = \{x_i, y_i / i = 1, 3, \dots, 2a+1\}$$

**Theorem 4.1:** If  $L(a)$  is the chemical graph (see Figure 2) then the  $K_1$  index of the graph  $L(a)$  is  $K_1(L(a)) = 4(a-1)$ , where  $a$  is a positive integer.

**Proof:** Consider

$$K(L(a)) = \sum_{e=\{x_i, x_{i+1}\} \in E_1} W(e) + \sum_{e=\{y_i, y_{i+1}\} \in E_2} W(e) + \sum_{e=\{x_i, y_i\} \in E_3} W(e)$$

Here

$$\sum_{e=\{x_i, x_{i+1}\} \in E_1} W(e) = (2a-2)1 \\ \text{(since there are } 2a-2 \text{ edges with } |i-j|=1 \text{ in } E_1). \\ \sum_{e=\{y_i, y_{i+1}\} \in E_2} W(e) = (2a-2)1 \\ \text{(since there are } 2a-2 \text{ edges with } |i-j|=1 \text{ in } E_2).$$

$$\sum_{e=\{x_i, y_i\} \in E_3} W(e) = 0 \text{ (since for every edge in } E_3, |i-j|=0). \\ \text{Therefore } K(L(a)) = 2a-2 + 2a-2 = 4a-4 = 4(a-1).$$

**Theorem 4.2:** If  $L(a)$  is the chemical graph (see Figure 2) then the  $K_1$  index of the graph  $L(a)$  is  $K_1(L(a)) = 2\sqrt{a-1}$ , where  $a$  is a positive integer.

**Proof:** Consider

$$K_1(L(a)) = \sum_{e=\{x_i, x_{i+1}\} \in E_1} W(e) + \sum_{e=\{y_i, y_{i+1}\} \in E_2} W(e) + \sum_{e=\{x_i, y_i\} \in E_3} W(e)$$

here

$$\sum_{e=\{x_i, x_{i+1}\} \in E_1} W(e) = (2a-2)1 \\ \text{(since there are } 2a-2 \text{ edges with } |i-j|=1 \text{ in } E_1).$$

$$\sum_{e=\{y_i, y_{i+1}\} \in E_2} W(e) = (2a-2)1 \\ \text{(since there are } 2a-2 \text{ edges with } |i-j|=1 \text{ in } E_2).$$

$$\sum_{e=\{x_i, y_i\} \in E_3} W(e) = 0 \\ \text{(since for every edge in } E_3, |i-j|=0).$$

Therefore

$$K_1(L(a)) = \sqrt{2a-2} + \sqrt{2a-2} = \sqrt{4a-4} = 2\sqrt{a-1}$$

**Theorem 4.3:** If  $L(a)$  is the chemical graph (see Figure 2) then the  $K_2$  index of the graph  $L(a)$  is  $K_2(L(a)) = 26a-2$ , where  $a$  is a positive integer.

**Proof:** In the set A two vertices namely  $x_1$  and  $x_{2a+1}$  having weight 4, two vertices namely  $x_2$  and  $x_{2a}$  having weight 5,  $(a-2)$  vertices namely  $x_4, x_6, x_8, \dots, x_{2a-2}$  having weight 6,  $(a-1)$  vertices namely  $x_3, x_5, x_7, \dots, x_{2a-1}$  having weight 7.

In the set B two vertices namely  $y_1$  and  $y_{2a+1}$  having weight 4, two vertices namely  $y_2$  and  $y_{2a}$  having weight 5,  $(a-2)$  vertices namely  $y_4, y_6, y_8, \dots, y_{2a-2}$  having weight 6,  $(a-1)$  vertices namely  $y_3, y_5, y_7, \dots, y_{2a-1}$  having weight 7.

Therefore

$$K_2(L(a)) = \sum_{\{x_i\} \in A} W(x_i) + \sum_{\{y_i\} \in B} W(y_i)$$

Here

$$\sum_{\{x_i\} \in A} W(x_i) = 2(4) + 2(5) + (a-2)6 + (a-1)7 \\ \sum_{\{y_i\} \in B} W(y_i) = 2(4) + 2(5) + (a-2)6 + (a-1)7$$

Therefore

$$K_2(L(a)) = 2(4) + 2(5) + (a-2)6 + (a-1)7 + 2(4) + 2(5) + (a-2)6 + (a-1)7 = 26a-2$$

**Theorem 4.4:** If  $L(a)$  is the chemical graph (see Figure 2) then the  $K_3$  index of the graph  $L(a)$  is  $K_3(L(a)) = 24$  when  $a = 1 = 42a-20$ , where  $a > 1$  is a positive integer.

**Proof:** In the set A two vertices namely  $x_1$  and  $x_{2a+1}$  having weight 4, two vertices namely  $x_2$  and  $x_{2a}$  having weight 6, (a-2) vertices namely  $x_4, x_6, x_8, \dots, x_{2a-2}$  having weight 9, (a-1) vertices namely  $x_3, x_5, x_7, \dots, x_{2a-1}$  having weight 12.

In the set B two vertices namely  $y_1$  and  $y_{2a+1}$  having weight 4, two vertices namely  $y_2$  and  $y_{2a}$  having weight 6, (a-2) vertices namely  $y_4, y_6, y_8, \dots, y_{2a-2}$  having weight 9, (a-1) vertices namely  $y_3, y_5, y_7, \dots, y_{2a-1}$  having weight 12.

Therefore

$$K_3(L(a)) = \sum_{\{x_i\} \in A} W(x_i) + \sum_{\{y_i\} \in B} W(y_i)$$

Here

$$\sum_{\{x_i\} \in A} W(x_i) = 2(4) + 2(6) + (a-2)9 + (a-1)12$$

$$\sum_{\{y_i\} \in B} W(y_i) = 2(4) + 2(6) + (a-2)9 + (a-1)12$$

Therefore

$$K_3(L(a)) = 2(4) + 2(6) + (a-2)9 + (a-1)12 + 2(4) + 2(6) + (a-2)9 + (a-1)12 = 42a - 20.$$

**Theorem 4.5:** If  $L(a)$  is the chemical graph (see Figure 2) then the  $K_4$  index of the graph  $L(a)$  is  $K_4(L(a)) = 18$  when  $a = 1 = 6a^2 + 8a + 2$ , where  $a > 1$  is a positive integer.

**Proof:** We know  $K_4(G) = \sum_{v \in V(G)} W(v)$

$$K_4(L(a)) = \sum_{\{x_i\} \in A} W(x_i) + \sum_{\{y_i\} \in B} W(y_i) \text{ Here}$$

$$\sum_{\{x_i\} \in A} W(x_i) = 2[(2a+1) + (2a) + (2a-1) + \dots + (2a-(a-2))] + (2a-(a-1))$$

$$\sum_{\{y_i\} \in B} W(y_i) = 2[(2a+1) + (2a) + (2a-1) + \dots + (2a-(a-2))] + (2a-(a-1))$$

Therefore

$$K_4(L(a)) = 2[(2a+1) + (2a) + (2a-1) + \dots + (2a-(a-2))] + (2a-(a-1)) + 2[(2a+1) + (2a) + (2a-1) + \dots + (2a-(a-2))] + (2a-(a-1))$$

$$= 2[2(a)(2a) - 2(2+3+\dots+(a-2))] + 2(a+1)$$

$$= 2[4a^2 - (a-2)(a-1) + 2] + (2a+2)$$

$$= 6a^2 + 8a + 2.$$

**Remark 4.6:**  $K_1, K_2, K_3$  can be represented in terms of  $K$  as follows:

$$K_1 = \sqrt{K}, K_2 = \frac{26}{4}K + 24, K_3 = \frac{42}{4}K + 22 \text{ (where } a \neq 1 \text{ in } K_3)$$

## COMPARISON OF RESULTS

In this section we compute the correlation of  $\log p$  with  $K, K_1, K_3$  and  $K_4$  considering the first 20 compounds of polyacenes ( $L(a)$ ,  $a=1$  to 20)

The values of  $K, K_1, K_2, K_3, K_4$  and  $\log p$  for these compounds are tabulated below (Table 1).

Table 1.

No	K	$K_1$	$K_2$	$K_3$	$K_4$	$\log p$
1	0	0	24	24	18	2.202
2	4	2	50	64	42	3.396
3	8	2.828	76	106	80	4.590
4	12	3.464	102	148	13	5.784
5	16	4	128	190	192	6.978
6	20	4.472	154	232	266	8.172
7	24	4.8989	180	274	352	9.366
8	28	5.2915	206	316	450	10.560
9	32	5.657	232	358	560	11.754
10	36	6	258	400	682	12.948
11	40	6.3245	284	442	816	14.142
12	44	6.6332	310	484	962	15.336
13	48	6.928	336	526	1120	16.530
14	52	7.211	362	568	1290	17.724
15	56	7.4833	388	610	1472	18.918
16	60	7.7459	414	652	1666	20.112
17	64	8	440	694	1872	21.306
18	68	8.2642	466	736	2090	22.500
19	72	8.4852	492	778	2320	23.694
20	76	8.7178	518	820	2562	24.888

Following are correlation coefficients of  $\log p$  with  $K, K_1, K_3$  and  $K_4$

Table 2.

r	$\log p$
$K$	0.99999997250969
$K_1$	0.96545941249478
$K_3$	0.99999854115530
$K_4$	0.97445287313308

## CONCLUSIONS

From the table it is clear that  $K$  is highly correlated with  $\log p$  compared to other indices. In [10] we observed that among  $W, PI, Sz, Sh$  and  $Fr$  indices of polyacenes,  $Sh_7$  is highly correlated with  $\log p$  with the value of  $r = 0.9996$ . However our new index  $K$  has better correlation 0.9999997250969 (almost one) than  $Sh$  index and this is more suitable for QSAR/QSPR studies. The regression equation between  $\log p$  and  $K$  is  $\log p = 2.20268571428571 + 0.29847142857143 K$  and the graphs of  $\log p$  and predicted  $\log p$  are given below.

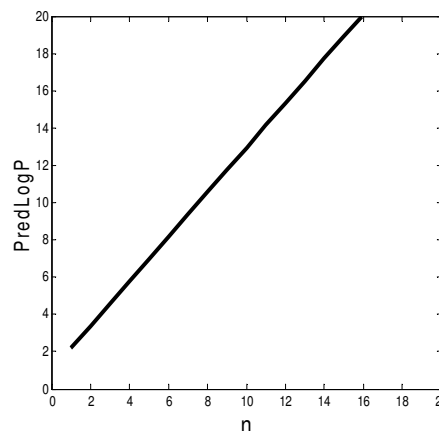


Fig 3 (a). Graph of predicted  $\log p$

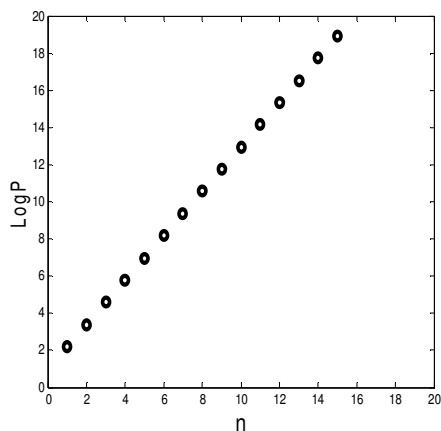


Fig 3 (b). Graph of log p

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