

Omega Polynomial in Polybenzene Multi Tori

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Abstract. The polybenzene units BTX₄₈, X=A (armchair) and X=Z (zig-zag) dimerize forming “eclipsed” isomers, the oligomers of which form structures of five-fold symmetry, called multi-tori. Multi-tori can be designed by appropriate map operations. The genus of multi-tori was calculated from the number of tetrapodal units they consist. A description, in terms of Omega polynomial, of the two linearly periodic BTX-networks was also presented.

Keywords: Polybenzene, Multi torus, Genus of structure, Linear periodic network, Omega polynomial.

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1. INTRODUCTION

The polybenzene unit BTA₄₈ (Figure 1, top, left) was shown [35] to dimerize to three different dimers BTA_{2_88}, BTA_{2_84} and BTA_{2_90}, by identifying the rings R(8) and R(12), respectively. Among these, the “eclipsed” dimer BTA_{2_90}, shows suitable angles to form a hyper-pentagon (Figure 1, bottom, left) structures of five-fold symmetry, eventually called multi-tori. The unit BTZ₂₄ (Figure 1, top, right) can form only an “eclipsed” dimer BTZ_{2_48}

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which of course can form a hyper-pentagon (Figure 1, bottom, right) and next multi-tori.

Multi-tori are complex structures consisting of more than one single torus [4,13,14]. They include negatively curved substructures [24,25,37], termed schwarzites, in the honor of H. A. Schwarz [29,30], who firstly investigated the differential geometry of this kind of surfaces. Multi tori can appear as self-assembly products of some repeating units/monomers, formed by spanning of cages/fullerenes, as in the case of spongy carbon or in natural zeolites. Multi tori MT can grow by a linear periodicity or by forming spherical arrays of various complexity [13]. They can be designed by appropriate map operations [3,5,11,15,33], as implemented in our original software CVNET [34] and Nano Studio [27].

The name of multi tori, bearing the benzene patch, will have B as a prefix. Next, because the opening faces show either “zig-zag” or “armchair” endings, “Z” or “A” will be added as a suffix to their name, as in BTZ or BTA. The number of repeating units and/or number of atoms will be added after the letters.

The design of simple units used to build up multi-tori was made by using some operations on maps, applied on the Tetrahedron T (see the letter “T” in the name of these units).

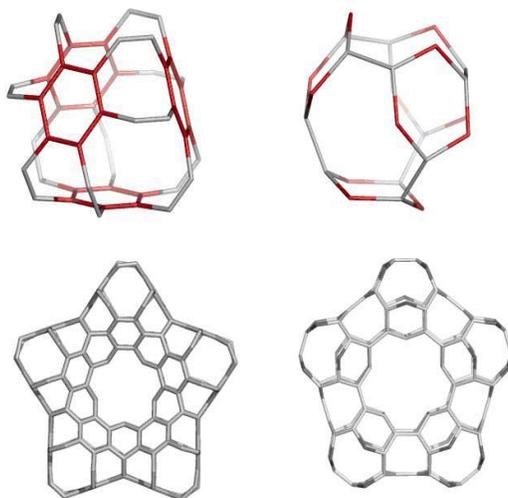


FIGURE 1. BTA₄₈ and BTZ₂₄ units and their corresponding hyper-pentagons BTX_{C_y5}.

2. DESIGN OF MULTI TORI

The hyper-ring BTX_{Cy5} , ($X=A, Z$, Figure 1, bottom), can self-arrange to a spherical multi torus $\text{BTX}20$ (Figure 2, left column), of genus $g=21$, with a well-defined core: $-f_5(\text{Le}_{2,2}(\text{Do}))=\text{core}(\text{BMTA}20)_180$, while $-d_5(\text{S}_2(\text{Ico}))=\text{core}(\text{BMTZ}20)_120$. In the above, $-f_5$ means deletion of all pentagonal faces in the transformed by Leapfrog (2,2) of the Dodecahedron Do , and d_5 is deletion of vertices of degree $d=5$, in the transform of Icosahedron= Ico by the septupling S_2 operation. Also, $-d_5(\text{S}_2(\text{Ico}))=\text{Op}(\text{Le}(\text{Ico}))$.

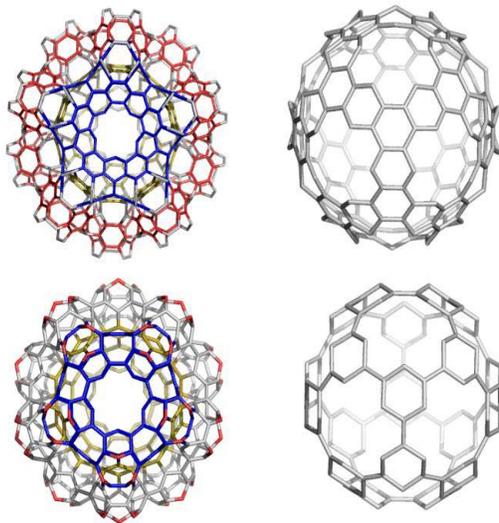


FIGURE 2. Bottom row: multi torus **BT**A20_1_780 (left) and its core_180 (right) designed by $-f_5(\text{Le}_{2,2}(\text{Do}))$. Top row: multi torus **BT**Z20_1_480 (left) and its core_120 (right) designed by $-d_5(\text{S}_2(\text{Ico}))$.

A linear array of $\text{BTX}20$, with the repeating unit formed by two units superimposing one pentagonal hyper-face (i.e., BTX_{Cy5}), rotated to each other by an angle of $\text{PI}/5$ as in the “dimer” $\text{BTX}20_2$ ($X=A$, Figure 3, left). Next, the structure can evolve with a one-dimensional periodicity, as shown in $\text{BTX}20_4$ (Figure 3, right).

The number u of tetrahedral units BTX in the linear array of $\text{BTX}20_k$ (Table 1) is $u=20k-5(k-1)=15k+5$, according to the construction mode. The term $-5(k-1)$ accounts for the superimposed hyper-rings BTX_{Cy5} , k being the number of units $\text{BTX}20$.

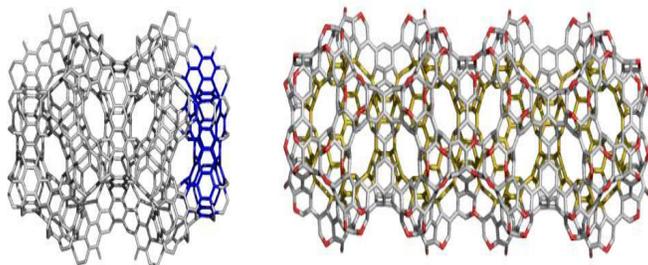


FIGURE 3. The repeating unit BTA20_2_1350 (left) and a rod-like array BTZ20_4_1560 (right).

The number u is also related to the number of faces as: $u = f_s/6$ in case BTA and $u = f_6/4$ in case BTZ (see Table 1).

The genus g of the surface where a structural graph is embedded counts the number of simple tori consisting that graph [20].

Theorem. [16,36] *In multi tori built up from open tetrahedral units, the genus of structure equals the number of its units plus one, irrespective of the unit tessellation.*

Demonstration comes out from construction: there are five tetrapodal units to be inserted into exactly five simple tori and all-together are joined to the central torus (see Figure 1, bottom), thus demonstrating the first part of the theorem.

For the second part, we apply the Euler's theorem [18]: $v - e + f = 2(1 - g)$, where $v = |V(G)|$ is the number of vertices/atoms, $e = |E(G)|$, the number of edges/bonds and f is the number of faces of the graph/molecule. Data in Table 1 provide the values of g in several BTX multi tori, tessellation differing as $X=A$ or Z , thus completing the demonstration

Table 1. Euler formula calculation in multi tori BTX.

	BTX	v	e	f_6	f_s	f_{tot}	$2(1-g)$	g	u	u -formula
1	BTACy5	210	285	35	30	65	-10	6	5	$f_s/6$
2	BTZCy5	120	165	20	15	35	-10	6	5	$f_6/4$
3	BTA20_1	780	1110	170	120	290	-40	21	20	$f_s/6$
4	BTZ20_1	480	690	80	90	170	-40	21	20	$f_6/4$
5	BTA20_5	3060	4410	710	480	1190	-160	81	80	$f_s/6$
6	BTZ20_5	1920	2790	320	390	710	-160	81	80	$f_6/4$

3. OMEGA POLYNOMIAL IN **Linear Multi Tori BTX20_k**

In a connected graph $G(V,E)$, with the vertex set $V(G)$ and edge set $E(G)$, two edges $e = uv$ and $f = xy$ of G are called *codistant e co f* if they obey the relation [22]:

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

which is reflexive, that is, $e \text{ co } e$ holds for any edge e of G , and symmetric, if $e \text{ co } f$ then $f \text{ co } e$. In general, relation co is not transitive; if “ co ” is also transitive, thus it is an equivalence relation, then G is called a *co-graph* and the set of edges $C(e) := \{f \in E(G); f \text{ co } e\}$ is called an *orthogonal cut oc* of G , $E(G)$ being the union of disjoint orthogonal cuts: $E(G) = C_1 \cup C_2 \cup \dots \cup C_k$, $C_i \cap C_j = \emptyset$, $i \neq j$. Klavzar [23] has shown that relation co is a theta Djokovi-Winkler relation [17,39].

We say that edges e and f of a plane graph G are in relation *opposite, e op f*, if they are opposite edges of an inner face of G . Note that the relation co is defined in the whole graph while op is defined only in faces. Using the relation op we can partition the edge set of G into *opposite edge strips, ops*. An *ops* is a quasi-orthogonal cut *qoc*, since ops is not transitive.

Let G be a connected graph and S_1, S_2, \dots, S_k be the *ops* strips of G . Then the *ops* strips form a partition of $E(G)$. The length of *ops* is taken as maximum. It depends on the size of the maximum fold face/ring F_{max}/R_{max} considered, so that any result on Omega polynomial will have this specification.

Denote by $m(G,s)$ the number of *ops* of length s and define the Omega polynomial as [1,6-10,12,28,38]:

$$\Omega(G, x) = \sum_s m(G, s) \cdot x^s \quad (2)$$

Its first derivative (in $x=1$) equals the number of edges in the graph:

$$\Omega'(G, 1) = \sum_s m(G, s) \cdot s = e = |E(G)| \quad (3)$$

On Omega polynomial, the Cluj-Ilmenau index [26], $CI=CI(G)$, was defined:

$$CI(G) = \{[\Omega'(G, 1)]^2 - [\Omega'(G, 1) + \Omega''(G, 1)]\} \quad (4)$$

Formulas to calculate Omega polynomial and CI index in the two infinite networks BTA20k and BTZ20k, designed on the ground of BTA₄₈ and BTZ₂₄ units, are presented in Tables 2 and 3. Formulas were derived from the numerical data calculated on rods consisting of k units BTX20. Omega polynomial was calculated at $R_{max}=R(8)$; examples are given in view of an easy verification of the general formulas. Formulas for the number of atoms, edges and rings (R_6 , R_8 and R_{15} , the last one being the simple ring of the hyper-ring BTX_{Cy5}), are included in Tables 2 and 3. Note the Omega polynomial description is an alternative to the crystallographic description.

Table 2. Formulas for Omega polynomial and net parameters in linear periodic BTA20_k network.

BMTA20 _k	$R_{max}(8);$ $\Omega(BMTA20_k-R_8) = 10(k+2)X^3 + 5(k-1)X^4 + (11k+1)X^5 + 20(k+3)X^8 + 10(k-1)X^{10} + 15(k-1)X^{12} + (11k+1)X^{20} + 10X^{2(3k+1)}$ $\Omega'(1) = 825k + 285 = E(G) = edges;$ $CI(G) = 15(45351k^2 + 30715k + 5332);$ $atoms = 10(57k + 21) = V(G);$ $R_6 = 5(27k + 7); R_8 = 30(3k + 1); R_{15} = 11k + 1;$ $u_{48} = 20k - 5(k - 1) = 5(3k + 1) = R_8/6;$ $g = 1 + u_{48}$
Examples	$k=5;$ $CI=19390230; atoms=3060; edges=4410; R_6=710; R_8=480; R_{15}=56; u_{48}=80; g=81.$ $k=6;$ $CI=27333870; atoms=3630; edges=5235; R_6=845; R_8=570; R_{15}=67; u_{48}=95; g=96.$

Table 3. Formulas for Omega polynomial and net parameters in linear periodic BTZ20_k network.

BMTZ20 _k	$R_{max}(8)$ $\Omega(BMTZ20_k-R_8) = 10(k+2)X^2 + 30kX^3 + (11k+1)X^5 + 10(k+5)X^6 + 10(k-1)X^8 + 10(k-1)X^{10} + 6kX^{20}$ $\Omega'(1) = 525k + 165 = E(G) = edges$ $CI(G) = 5(55125k^2 + 33653k + 5392)$ $atoms = 120(3k + 1) = V(G) = 24u_{24} = 6R_6$ $R_6 = 20(3k + 1) = V(G) /6; R_8 = 15(5k + 1); R_{15} = 11k + 1$ $u_{24} = 20k - 5(k - 1) = 5(3k + 1) = R_6/4;$ $g = 1 + u_{24}$
Examples	$k=5; 70X^2+150X^3+56X^5+100X^6+40X^8+40X^{10}+30X^{20}$ $CI=7758910; atoms=1920; edges=2790; R_6= 320; R_8=390; R_{15}=56; u_{24}=80; g=81.$ $k=6; 80X^2+180X^3+67X^5+110X^6+50X^8+50X^{10}+36X^{20}$ $CI=10959050; atoms=2280; edges=3315; R_6=380; R_8=465; R_{15}=67; u_{24}=95; g=96.$

4. CONCLUSIONS

Polybenzene units BTX₄₈ was shown to dimerize forming “eclipsed” isomers, the oligomers of which form structures of five-fold symmetry, called multi-tori.

Multi-tori can grow by a linear periodicity or by forming spherical arrays of various complexity [2]. They can be designed by appropriate map operations [10-14], as implemented in the software CVNET [15] and Nano Studio [16] developed at TOPO Group Cluj. The genus of multi-tori was calculated from the number of tetrapodal units they consist. A description, in terms of Omega polynomial, of the two linear BTX-networks was also presented. We mention that in the last years several authors have published articles dealing with the calculation of various topological indices [2,21,26,32] and counting polynomials [19,31].

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REFERENCES

- [1] A. R. Ashrafi, M. Jalali, M. Ghorbani and M. V. Diudea, Computing PI and Omega polynomials of an infinite family of fullerenes, *MATCH Commun. Math. Comput. Chem.*, **60**, (2008), 905-916.
- [2] Z. Bagheri , A. Mahmiani , O. Khormali, Edge-Szeged and vertex-PI indices of some benzenoid systems, *Iranian Journal of Mathematical Sciences and Informatics*, **3** (1), (2008), 31-39.
- [3] M. V. Diudea, Covering forms in nanostructures, *Forma (Tokyo)*, **19**, (2004), 131-163.
- [4] M. V. Diudea, *Nanostructures, Novel Architecture*, Nova, N.Y., 2005.
- [5] M. V. Diudea, Nanoporous carbon allotropes by septupling map operations, *J. Chem. Inf. Model.*, **45**, (2005), 1002-1009.
- [6] M. V. Diudea, Omega Polynomial, *Carpath. J. Math.*, **22**, (2006), 43-47.
- [7] M. V. Diudea, Counting polynomials in tori $T(4,4)S[c,n]$, *Acta Chim. Slov.*, **57**, (2010), 551-558.
- [8] M. V. Diudea, S. Cigher and P. E. John, Omega and Related Counting Polynomials, *MATCH Commun. Math. Comput. Chem.*, **60**, (2008), 237-250.
- [9] M. V. Diudea, S. Cigher, A. E. Vizitiu, M. S. Florescu and P. E. John, Omega polynomial and its use in nanostructures description, *J. Math. Chem.*, **45**, (2009), 316-329.
- [10] M. V. Diudea, S. Cigher, A. E. Vizitiu, O. Ursu and P. E. John, Omega Polynomial in Tubular Nanostructures, *Croat. Chem. Acta*, **79**, (2006), 445-448.
- [11] M. V. Diudea, P. E. John, A. Graovac, M. Primorac and T. Pisanski, Leapfrog and related operations on toroidal fullerenes, *Croat. Chem. Acta*, **76**, (2003), 153-159.
- [12] M. V. Diudea, S. Klavzar, Omega polynomial revisited, *Acta Chim. Slov.*, **57**, (2010), 565-570.
- [13] M. V. Diudea and Cs. L. Nagy, *Periodic Nanostructures*, Springer, 2007.
- [14] M. V. Diudea and M. Petitjean, Symmetry in multi tori, *Symmetry, Culture, Sci.*, **19** (4), (2008), 285-305.
- [15] M. V. Diudea, M. Stefu, P. E. John and A. Graovac, Generalized operations on maps, *Croat. Chem. Acta*, **79**, (2006), 355-362.
- [16] M. V. Diudea and B. Szefer, Nanotube junctions and the genus of multi-tori, *Phys. Chem. Chem. Phys.*, **14**, (2012), 8111-8115.
- [17] D. Z. Djokovic, Distance preserving subgraphs of hypercubes, *J. Combin. Theory Ser. B*, **14**, (1973), 263-267.
- [18] L. Euler, Elementa doctrinae solidorum, *Novi Comm. Acad. Scient. Imp. Petrop.*, **4**, (1758), 109-160.
- [19] G. Fath-Tabar, A. R. Ashrafi, The Hyper-Wiener polynomial of graphs, *Iranian Journal of Mathematical Sciences and Informatics*, **6** (2), (2011), 67-74.
- [20] F. Harary , *Graph Theory*, Addison - Wesley, Reading, M.A., 1969.
- [21] A. Iranmanesh , Y. Alizadeh , Computing Wiener index of HAC5C7[p,q] nanotubes by GAP program, *Iranian Journal of Mathematical Sciences and Informatics*, **3** (1), (2008), 1-12.
- [22] P. E. John, A. E. Vizitiu, S. Cigher and M. V. Diudea, CI Index in Tubular Nanostructures, *MATCH Commun. Math. Comput. Chem.*, **57** (2), (2007), 479-484.
- [23] S. Klavzar, Some comments on co graphs and CI index, *MATCH; Commun. Math. Comput. Chem.*, **59**, (2008), 217-222.
- [24] T. Lenosky, X. Gonze, M. Teter and V. Elser, Energetics of Negatively Curved. Graphitic Carbon, *Nature*, **355**, (1992), 333-335.
- [25] A. L. Mackay and H. Terrones, Diamond from Graphite, *Nature*, **352**, (1991), 762-762.

- [26] A. Mahmiani , O. Khormali , A. Iranmanesh, The explicit relation among the edge versions of detour index, *Iranian Journal of Mathematical Sciences and Informatics*, **3** (2), (2008), 1-12.
- [27] Cs. L. Nagy and M. V. Diudea, *Nano Studio* software, Babes-Bolyai University, 2009.
- [28] M. Saheli, M. Neamati, K. Nagy and M. V. Diudea, Omega polynomial in Sucor network. *Studia Univ. Babes-Bolyai*, **55** (1), (2010), 83-90.
- [29] H. A. Schwarz, Uber Minimalflachen, *Monatsber*, Berlin Akad., 1865.
- [30] H. A. Schwarz, *Gesammelte Matematische Abhandlungen*, Springer, Berlin, 1890.
- [31] S. Sedghi , N. Shobe , M. A. Salahshoor, The polynomials of a graph, *Iranian Journal of Mathematical Sciences and Informatics*, **3** (2), (2008), 55-67.
- [32] A. Sousaraei , A. Mahmiani , O. Khormali, Vertex-PI index of some nanotubes, *Iranian Journal of Mathematical Sciences and Informatics*, **3** (1), (2008), 49-62.
- [33] M. Stefu, M. V. Diudea and P. E. John, Composite operations on maps, *Studia Univ. "Babes-Bolyai", Chemia*, **50** (2), (2005), 165-174.
- [34] M. Stefu and M. V. Diudea, *CageVersatile* 1.5 software, Babes-Bolyai University, 2005.
- [35] B. Szeffler and M. V. Diudea, Polybenzene revisited, *Acta Chim. Slo.*, 2012 (submitted).
- [36] B. Szeffler and M. V. Diudea, Polybenzene multi tori, *Central Eur. J.Chem.*, 2012 (submitted).
- [37] H. Terrones and A. L. Mackay, Triply periodic minimal surfaces decorated with curved graphite, *Chem. Phys. Lett.*, **207**, (1993), 45-50.
- [38] A. E. Vizitiu, S. Cigher, M. V. Diudea and M. S. Florescu, Omega Polynomial in ((4,8)3) tubular nanostructures, *MATCH Commun. Math. Comput. Chem.*, **57** (2), (2007), 457-462.
- [39] P. M. Winkler, Isometric embedding in products of complete graphs, *Discrete Appl. Math.*, **8**, (1984), 209-212.