

Proposition 5 (ii) is deduced from

$$A_i = A_{R_1} + A_{R_2} \quad \text{and} \quad A_0 = A_{R_1} \cap A_{R_2},$$

which are verified as for (i).

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*Acta Cryst.* (1976). **A32**, 65

## Univalent (Monodentate) Substitution on Convex Polyhedra. II. Listing of Cycle Indices

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(Received 14 July 1975; accepted 15 July 1975)

To extend the usefulness of the tabulation of the numbers  $N$  of positional isomers [Knop, Barker & White (1975). *Acta Cryst.* **A31**, 461–472], all the distinct cycle-index polynomials  $Z$  on which the tabulation is based have been listed in a convenient form. This condensed summary facilitates identification of  $Z$ -isomorphisms; in turn,  $N$  for univalent substitution on many polyhedra not listed previously can be evaluated simply by reference to the existing tabulation.

In part I (Knop, Barker & White, 1975) we presented the numbers  $N$  of distinct (up to rotation) positional isomers obtained by univalent substitution at the vertices of convex polyhedra; only structureless substituents were considered. The tabulation is extensive, but naturally it cannot include *all* non-isomorphic polyhedra even for small numbers of vertices  $V$ . A user of the tables wishing to evaluate  $N$  for polyhedra not listed in Table 5 of part I would not only have to determine the appropriate cycle indices  $Z$ , but he would have to compute the coefficients of the expanded cycle-index polynomials (*i.e.* the values of  $N$ ) for the compositions of interest, a tedious task. However, owing to cycle-index isomorphism the number of distinct  $Z$  polynomials involved in the tabulations of part I is not unduly large, and there is a good chance that the set of  $N$  to be determined already appears there under a different but  $Z$ -isomorphic polyhedron, which makes fresh computation unnecessary.

To facilitate identification of additional  $Z$ -isomorphisms, over and above those *specifically* listed in part I, a table of *all* the cycle indices on which the tabulation of part I is based, has been compiled.

Considerable space is saved by introducing the following notation. An  $s$ -product  $s_a^u s_b^v$  will be represented as  $a, u * b, v$ . Each  $s$ -product occurring in the cycle indices

for a particular value of  $V$  will be denoted by a capital letter (Table 1). The highest-order term  $s_1^V$  (represented by  $A$ ) is always present,† and so further space is saved by omitting  $A$  from the letter symbol of  $Z$ . For example, the  $Z$  of a tetrahedron 4-2 of symmetry  $T_d$ ,

$$\frac{1}{24}(s_1^4 + 6s_1^3 s_2 + 8s_1^2 s_3 + 6s_1 s_2^2 + 3s_2^3),$$

is represented by 6B8C6D3E. The  $s$ -products denoted by the letters are found in Table 1 under  $V=4$ . The sum of the coefficients associated with the letters, including the coefficient of  $A$ , which is always unity, is equal to the divisor  $p(\mathbf{G})$ , in this case 24.

For economy of space, the table of cycle indices (Table 2) is arranged as follows. In the first part (pp. 3–9)  $Z$  polynomials occurring in only a few cases are listed in the order of increasing  $V$ . The second part (pp. 9–16) contains cycle indices having large numbers of terms and those involved in considerable numbers of  $Z$ -isomorphic representations.

† The term  $A$  by itself represents  $Z(4mh)$  of the corresponding polygon of  $V$  vertices (*cf.* part I).

‡ Table 2 has been deposited with the British Library Lending Division as Supplementary Publication No. SUP 31246 (16 pp., 1 microfiche). Copies may be obtained through The Executive Secretary, International Union of Crystallography, 13 White Friars, Chester CH1 1NZ, England.

Table 1. *Cycle-index terms*

<i>V</i>	<i>B</i>	<i>C</i>	<i>D</i>	<i>E</i>	<i>F</i>	<i>G</i>	<i>H</i>	<i>J</i>
3	3,1	1,1*2,1						
4	4,1	1,1*3,1	1,2*2,1	2,2				
5	5,1	1,1*4,1	2,1*3,1	1,2*3,1	1,1*2,2	1,3*2,1		
6	6,1	1,1*5,1	2,1*4,1	1,2*4,1	3,2	1,2*2,2	1,4*2,1	2,3
7	7,1	1,1*6,1	2,1*5,1	1,2*5,1	1,1*3,2	1,1*2,3	1,3*2,2	1,5*2,1
8	8,1	1,1*7,1	2,1*6,1	1,2*6,1	2,1*3,2	4,2	1,2*3,2	1,2*2,3
9	9,1	1,1*8,1	2,1*7,1	1,2*7,1	3,1*6,1	1,1*4,2	3,3	1,1*2,4
10	10,1	1,1*9,1	2,1*8,1	1,2*8,1	2,1*4,2	5,2	1,2*4,2	1,1*3,3
11	11,1	1,1*10,1	2,1*9,1	1,2*9,1	2,1*3,1*6,1	1,1*5,2	2,1*3,3	1,2*3,3
12	12,1	1,1*11,1	2,1*10,1	1,2*10,1	2,1*5,2	6,2	1,2*5,2	4,3
13	13,1	1,1*12,1	2,1*11,1	1,2*11,1	1,1*6,2	1,1*4,3	1,1*3,4	1,1*2,6
14	14,1	1,1*13,1	2,1*12,1	1,2*12,1	2,1*3,2*6,1	2,1*6,2	7,2	1,2*6,2
15	15,1	2,1*13,1	5,1*10,1	1,2*13,1	3,1*6,2	5,3	3,5	1,1*2,7
16	16,1	8,2	4,4	1,1*3,5	1,2*2,7	1,4*2,6	2,8	
17	17,1	1,1*2,8						
18	18,1	9,2	6,3	2,1*4,4	1,2*4,4	3,6	1,2*2,8	1,4*2,7
19	19,1	1,1*2,9						
20	10,2	2,1*6,3	5,4	4,5	1,2*3,6	1,2*2,9	1,4*2,8	1,6*2,7
24	12,2	8,3	6,4	4,6	3,8	1,2*2,11	1,4*2,10	1,8*2,8
26	2,1*6,4	2,1*4,6	1,2*4,6	1,2*3,8	1,2*2,12	1,8*2,9	2,13	
30	10,3	6,5	5,6	3,10	1,2*2,14	1,4*2,13	2,15	
32	2,1*10,3	2,1*6,5	1,2*5,6	1,2*3,10	1,8*2,12	2,16		
38	1,2*4,9	1,2*3,12	1,2*2,18	2,19				
48	6,8	4,12	3,16	2,24				
60	10,6	6,10	5,12	3,20	1,4*2,28	2,30		
62	2,1*10,6	2,1*6,10	1,2*5,12	1,2*3,20	1,2*2,30	1,12*2,25	2,31	
92	1,2*5,18	1,2*3,30	2,46					
120	10,12	6,20	5,24	3,40	2,60			
<i>K</i>	<i>L</i>	<i>M</i>	<i>N</i>	<i>P</i>	<i>Q</i>	<i>R</i>	<i>S</i>	<i>T</i>
8	1,4*2,2	1,6*2,1	2,4					
9	1,3*2,3	1,5*2,2	1,7*2,1					
10	1,2*2,4	1,4*2,3	1,6*2,2	1,8*2,1	2,5			
11	1,1*2,5	1,3*2,4	1,5*2,3	1,9*2,1				
12	3,4	1,2*2,5	1,4*2,4	1,6*2,3	1,10*2,1	2,6		
13	1,3*2,5	1,5*2,4	1,11*2,1					
14	2,1*4,3	1,2*4,3	2,1*3,4	1,2*3,4	1,2*2,6	1,4*2,5	1,6*2,4	1,12*2,1
15	1,3*2,6	1,5*2,5	1,13*2,1					2,7
18	1,8*2,5	2,9						
20	2,10							
24	2,12							

To aid photographic reduction the type used in Table 2 does not differentiate between the running number of the polyhedron (bold face type in Table 5 of part I), the running number of the point group (italics in Table 3 of part I), and the symbol representing the *Z* terms. Thus **4B2GM 8-1** (61,60,59) appears

in Table 2, without a possibility of confusion, as **4B2GM 8-1** (61,60,59).

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