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Correction

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CORRECTION

Molecular Simulation, 1991, Vol. 6, pp. 103-119

MONTE CARLO SIMULATION OF TETRAHEDRAL CHAINS II: PROPERTIES OF "FIRST SELF-AVOIDING WALKS" AND THEIR USABILITY AS STARTING CONFIGURATIONS FOR DYNAMIC RELAXATION MECHANISMS

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The shape-factors given in Table 1 to Table 3 actually have been obtained by averaging the ratios L_i^2/s^2 of individual chains and, therefore, are *not* ratios of two averaged quantities $\langle L_i^2 \rangle / \langle s^2 \rangle$ as stated in the paper.

Defining $sf'_i = L_i^2/s^2$ the labels sf_i in Table 1 to Table 3 should be changed to $\langle sf'_i \rangle$. In addition the values of the ratios $sf_i = \langle L_i^2 \rangle / \langle s^2 \rangle$ are given in the following Table.

	<i>NRRW</i>	<i>FSAW_L</i>	<i>FSAW_P</i>	<i>FSAW_R</i>	<i>SAW_R</i>	<i>SAW</i>
<i>N</i> = 50						
<i>sf</i> ₁	0.05928	0.06363	0.05569	0.05269	0.05323	0.05265
<i>sf</i> ₂	0.16926	0.17384	0.16533	0.16166	0.16225	0.16139
<i>sf</i> ₃	0.77147	0.76253	0.77898	0.78565	0.78452	0.78596
<i>N</i> = 500						
<i>sf</i> ₁	0.06285	0.06684	0.05434	0.05599	0.05595	0.05503
<i>sf</i> ₂	0.17207	0.17526	0.16267	0.16521	0.16585	0.16386
<i>sf</i> ₃	0.76508	0.75790	0.78299	0.77879	0.77819	0.78111
<i>N</i> = 1000						
<i>sf</i> ₁	0.06310	0.06647	0.05404	0.05602	0.05655	0.05529
<i>sf</i> ₂	0.17212	0.17404	0.16201	0.16542	0.16569	0.16384
<i>sf</i> ₃	0.76478	0.75949	0.78395	0.77856	0.77776	0.78087