

## Erratum: Random-Walk Model of Chain-Polymer Adsorption at a Surface

Robert J. Rubin

Citation: The Journal of Chemical Physics 44, 2222 (1966); doi: 10.1063/1.1727025

View online: http://dx.doi.org/10.1063/1.1727025

View Table of Contents: http://scitation.aip.org/content/aip/journal/jcp/44/5?ver=pdfcov

Published by the AIP Publishing

## Articles you may be interested in

Random-walk model of chain polymer adsorption behavior at critical energy and relation to a reflecting boundary condition

AIP Conf. Proc. 109, 73 (1984); 10.1063/1.34340

Ordered spans of unrestricted and self-avoiding random-walk models of polymer chains. I. Space-fixed axes

J. Chem. Phys. 63, 5362 (1975); 10.1063/1.431342

Theory of inhomogeneous polymers: Fundamentals of the Gaussian random-walk model

J. Chem. Phys. 62, 999 (1975); 10.1063/1.430517

Random-Walk Model of Adsorption of a Chain-Polymer Molecule on a Long Rigid-Rod Molecule

J. Chem. Phys. 44, 2130 (1966); 10.1063/1.1726991

Random-Walk Model of Chain-Polymer Adsorption at a Surface

J. Chem. Phys. 43, 2392 (1965); 10.1063/1.1697138



Although the theory makes no pretence of being anything but a rough guide, it is clear that most of the volume of ionization arises from the compression of the ions, and only a small part from the electrostriction of the medium.

## Erratum: Normal Vibrations of Polyacrylonitrile and Deuterated Polyacrylonitriles

[J. Chem. Phys. 41, 1233 (1964)] REIZO YAMADERA

Textile Research Institute, Toyo Spinning Company, Ltd. Katata, Shiga, Japan

 $\mathbf{I}^{\mathrm{N}}$  Eq. (2) on p. 1237, seventh term of the expression of 2V(j) should read

instead of

$$H(\operatorname{CCC}^{\prime\prime})\,{R_0}^{\prime 2} \sum_{\Theta_{\mathbf{III}},\Theta_{\mathbf{IV}}} \!\! \big[ \Delta\Theta(\,j) \,\big]^2.$$

In  $\alpha$  matrix on p. 1238, for third column in 17th row, read  $F_1$  instead of  $F_3$ , and for eighth column in 18th row, read U instead of  $L_0$ .

In  $\beta$  matrix on p. 1239, for 14th, 15th, and 16th columns in 10th row, read T,  $T_0$ , and  $P_2$  instead of  $T_0$ ,  $T_2$ , and P, respectively.  $\beta$  matrix should be transposed upon including the correction described above.

On p. 1240, for equations of  $S_2$  and  $T_1$ , read

$$S_2 = (5\rho_0 \rho - \rho_0^2) \mu_C/6$$

and

$$T_1 = (5\rho_0\rho - \rho_0\rho_1)\mu_C/6.$$

On p. 1242, the sign of vibrational mode on  $s_7$  in  $S(j)_{II_{\delta}}$  matrix should read  $\nu(CH)$ , not  $\delta(CH)$ .

## Erratum: Random-Walk Model of Chain-Polymer Adsorption at a Surface

[J. Chem. Phys. 43, 2392 (1965)]

ROBERT J. RUBIN

National Bureau of Standards, Washington, D. C.

THE leading minus sign in the third line of Eq. (39) should be moved up to the second line.

The heading of the middle column of Table II should be  $\nu(0, \infty)$ .