

Hypercubic lattice SAW exponents ν and γ : 3.99 dimensions revisited

Jack F Douglas[†], Takao Ishinabe[‡], Adolfo M Nemirovsky[§] and Karl F Freed^{||}

[†] Polymers Division, National Institute of Standards and Technology, Gaithersburg, MD 20899, USA

[‡] Faculty of Engineering, Yamagata University, Yonezawa 992, Japan

[§] Departamento de Estructura, y Constituyentes de la Materia, Universidad de Barcelona, Barcelona E-08028, Spain

^{||} James Franck Institute, University of Chicago, Chicago, IL 60637, USA

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Abstract. The self-avoiding walk (SAW) exponents ν and γ are computed over a range of dimensions ($1 \leq d < \infty$) from *exact* expressions for the mean-square end-to-end distance $\langle R_n^2 \rangle$ and the partition function Q_n of SAWs having a limited number of steps, $n \leq 11$. SAW exponents (ν, γ) for *arbitrary* dimension d are estimated by applying standard extrapolation techniques to our direct enumeration data which has been analytically continued to variable dimension. Exponent estimates obtained from continuum theories of self-avoiding paths are compared with the SAW calculations.

1. Introduction

It is well known that the geometrical properties of self-avoiding and random walks exhibit a strong dependence on spatial dimensionality d and that the 'universal' critical behaviour in many systems undergoing phase transitions is intimately related to the geometrical properties of these walks. Specifically, we mention Symanzik's formulation [1] of ϕ^4 field theory in terms of a 'gas' of interacting Brownian paths, Domb's calculation of $O(m)$ lattice spin model properties in terms of interacting SAWs [2] and De Gennes' discovery of an exact relation between the $m \rightarrow 0$ limit of the $O(m)$ model and SAWs [3]. The accurate characterization of the geometrical properties of SAWs is consequently a problem with many practical physical applications, besides the rather obvious applications to the solution properties of polymers [4, 5].

Many analytical and numerical studies of self-avoiding walks with nearest-neighbour interactions on a variety of lattices have appeared since the pioneering studies by Orr [6] and Fisher *et al* [7]. Recent works [8] often emphasize the calculation of the SAW 'critical indices' ν and γ , and the 'connectivity constant' μ . Rigorous results include a proof of the existence [9, 10] of μ and the relation [11], $\gamma = 2\nu = 1$, $d \geq 5$. Conformal invariance calculations in $d = 2$ suggest the exact results [12] $2\nu = \frac{3}{2}$ and $\gamma = \frac{43}{32}$.

The study of lattice SAW models has developed in parallel to analytic theories of self-avoiding paths based on $O(m \rightarrow 0)$ field theoretic methods [3] or direct formulations in terms of Wiener path-integration [4, 5, 13]. Application of the perturbative Wilson-Fisher ϵ -expansion method has provided SAW information that is complementary to lattice model studies [5, 13]. Simple dimensional analysis in the continuum theory of

interesting. Calculations are in progress for ν and γ for theta-point and neighbour-avoiding (NAWs) polymers ($\omega \rightarrow -\infty$) in variable dimensions [50].

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References

- [1] Symanzik K 1969 Euclidean quantum field theory *Local Quantum Field Theory* (New York: Academic) p 152. See also: Feynman R P 1953 *Phys. Rev.* **90** 1116; 1953 *Phys. Rev.* **91** 1291
Aizenman M 1982 *Commun. Math. Phys.* **86** 1
- [2] Domb C 1970 *J. Phys. C: Solid State Phys.* **3** 256; 1970 *J. Phys. C: Solid State Phys.* **5** 1399; 1972 *J. Phys. C: Solid State Phys.* **5** 1417
- [3] DeGennes P G 1972 *Phys. Lett.* **38A** 339
Fisher M E 1974 *Rev. Mod. Phys.* **46** 597
Wilson K G 1974 *Physica* **73** 119
- [4] DeGennes P G 1979 *Scaling Concepts in Polymer Physics* (Ithaca, NY: Cornell University Press)
- [5] Freed K F 1987 *Renormalization Group Theory of Macromolecules* (New York: Wiley) see [13]
- [6] Orr W J 1947 *Trans. Faraday Soc.* **43** 12
- [7] Fisher M E and Sykes M F 1959 *Phys. Rev.* **114** 45
Fisher M E and Hiley B J 1961 *J. Chem. Phys.* **34** 1253
- [8] Guttman A J 1989 *Phase Transitions* **13** 3
- [9] Hammersley J M and Morton K W 1954 *J. Roy. Stat. Soc. B* **16** 23
Hammersley J M 1957 *Proc. Camb. Phil. Soc.* **53** 642
- [10] Kesten H 1963 *J. Math. Phys.* **4** 960
- [11] Hara T and Slade G 1991 *Bull. AMS* **25** 417
Fröhlich J 1982 *Nucl. Phys. B* **200** 281
- [12] Nienhuis B 1982 *Phys. Rev. Lett.* **49** 1062; 1984 *J. Stat. Phys.* **34** 731
- [13] Jannink G and Des Cloizeaux J 1990 *J. Phys.: Condens. Matter* **2** 1
- [14] Douglas J F, Roovers J and Freed K F 1990 *Macromolecules* **23** 4168
- [15] Gaunt D S 1986 *J. Phys. A: Math. Gen.* **19** L149
Fisher M E and Gaunt D S 1964 *Phys. Rev. A* **133** 224
- [16] Gerber P R and Fisher M E 1975 *J. Chem. Phys.* **63** 4941
- [17] Elezovic S, Knežević M and Milošević S 1987 *J. Phys. A: Math. Gen.* **20** 1215
- [18] Milošević S and Živić I 1991 *J. Phys. A: Math. Gen.* **24** L833
- [19] Lam P M 1990 *J. Phys. A: Math. Gen.* **23** 831
Barat K, Karmakar S N and Chakrabarti B K 1991 *J. Phys. A: Math. Gen.* **24** 851
- [20] Meir Y and Brooks Harris A 1989 *Phys. Rev. Lett.* **25** 2819
- [21] LeGuillou J E and Zinn-Justin J 1987 *J. Physique* **48** 19
- [22] Nemirovsky A M, Freed K F, Ishinabe T and Douglas J F 1992 *J. Stat. Phys.* **67** 1083
- [23] Nemirovsky A M, Freed K F, Ishinabe T and Douglas J F 1992 *Phys. Lett.* **162A** 469
- [24] Abe R 1972 *Prog. Theor. Phys.* **47** 1200
- [25] Baker G A and Benofy L P 1982 *J. Stat. Phys.* **29** 699. See also: Griffiths R B and Gujrati 1983 *J. Stat. Phys.* **30** 563
- [26] Wegner F J 1972 *Phys. Rev. B* **5** 4529; 1972 *Phys. Rev. B* **6** 1891
- [27] Ishinabe T 1985 *J. Phys. A: Math. Gen.* **18** 3181; 1987 *J. Phys. A: Math. Gen.* **20** 6435
- [28] Barrett A J, Mansfield M and Benesch B C 1991 *Macromolecules* **24** 2615
- [29] LeGuillou J C and Zinn-Justin J 1977 *Phys. Rev. Lett.* **39** 95; 1980 *Phys. Rev. B* **21** 3976
- [30] Domb C 1969 *Adv. Chem. Phys.* **25** 229
- [31] Ishinabe T 1989 *Phys. Rev. B* **39** 9486