## Comment on "Interacting-Electron Model Exactly Solvable in Any Dimension"

In a recent Letter [1] Montorsi and Rasetti investigated a special fermionic lattice model

$$H = t \sum_{(ij)} \kappa_{ij} A_i^{\dagger} A_j + U \sum_{i} N_i D_i - \mu \sum_{i} (N_i + D_i), \qquad (1)$$

where the first term (kinetic energy  $H_0$ ) commutes with the second term (Hubbard interaction  $H_1$ ). Hence the double occupancy of lattice sites is conserved. Here  $A_i^{\dagger}$ ,  $A_i$ , with  $N_i = A_i^{\dagger} A_i$ , correspond to spinless fermions and  $D_i = D_i^2$  commutes with  $A_i$ . The amplitude  $\kappa_{ij} = 1 - D_i - D_j + \eta D_i D_j$  describes restricted hopping, where  $\eta$  is arbitrary. According to the authors they obtained the exact free energy density for (1) in any dimension d [Eq. (20) of Ref. [1]]; it has the general form

$$f = f_0 + N^{-1} \sum_{\mathbf{k}} F(\epsilon_{\mathbf{k}}) , \qquad (2)$$

where  $\epsilon_{\mathbf{k}}$  is the tight-binding dispersion.

In this Comment we would like to point out a discrepancy between the known properties of the model and the result of Ref. [1]. Using the notation  $A_i \equiv c_{i\uparrow}$ , etc.,  $N_i \equiv n_{i\uparrow}$  and  $D_i \equiv n_{i\downarrow}$ , and choosing  $\eta = 1$ , we see that for zero double occupancy the Hamiltonian (1) reduces

$$H = t \sum_{\langle ij \rangle} (1 - n_{i\downarrow}) c_{i\uparrow}^{\dagger} c_{j\uparrow} (1 - n_{j\downarrow})$$
(3)

which is identical with a simplified Hubbard model ("Falicov-Kimball model") at  $U=\infty$ , where only  $\uparrow$  spins are mobile. This model is known to be nontrivial even for  $U=\infty$  if n<1 [2]. By contrast, the dimension dependence of (2) is trivial, since the integrand of the k sum depends on momentum k only via  $\epsilon_k$ . Using  $N^{-1}\sum_{\bf k}F(\epsilon_{\bf k})=\int dE\,\rho_d(E)F(E)$ , where  $\rho_d(E)$  is the d-dimensional density of states of the noninteracting system, Eq. (2) is seen to have a mean-field form. The exact solution of (3), namely, in the limit of high dimensions  $d\to\infty$  [3], has recently been obtained by different methods [4,5]. In particular, in this limit and for  $U=\infty$  the free energy is given by [5]

$$f = \frac{1}{\pi} \int_{-\infty}^{\infty} d\omega f_T(\omega) \left[ \int_{-\infty}^{\infty} dE \, \rho_{\infty}(E) \operatorname{Im} \ln[\omega + \mu - E - \Sigma(\omega) + i0^+] + \operatorname{Im} \ln[1 + G(\omega)\Sigma(\omega)] \right] - \beta^{-1} \ln[1 + e^{\beta(\mu - \epsilon)}], \tag{4}$$

where

$$\epsilon = \frac{1}{\pi} \int d\omega f_T(\omega) \tan^{-1} \left[ \frac{-\operatorname{Im}[G^{-1}(\omega) + \Sigma(\omega)]}{\operatorname{Re}[G^{-1}(\omega) + \Sigma(\omega)]} \right]$$

and  $f_T(\omega)$  is the Fermi function. Here  $G(\omega)$  is the diagonal element of the Green function of mobile electrons and  $\Sigma(\omega)$  is the self-energy; they are determined by  $\delta f/\delta \Sigma = \delta f/\delta G = 0$ . To be able to compare (2) and (4) we introduce the limit  $d \to \infty$ , i.e.,  $\rho_\infty(E)$ , into (2) and put  $U = \infty$  [3]. The resulting expression is very different from (4). In view of this discrepancy in such a well-defined limit and the simplistic mean-field d dependence of (2) we conclude that (2) is not exact.

Furthermore, we would like to comment on the "metal-insulator transition" discussed in Ref. [1]. Taking, as in Ref. [1],  $\mathbf{k} \in \mathcal{D}$  in (2) ( $\mathcal{D}$ : half a Brillouin zone), half filling, and  $\rho = \pm 1$ , the ground-state energy in arbitrary dimensions d follows as

$$E_g = -2 \int_{U/4}^{2dt} dE \, \rho_d(E) (E - \frac{1}{4} U) \,. \tag{5}$$

Even if (5) were correct, the transition at U=8dt would not be of the Mott-Hubbard-type [6] (e.g., there is no band narrowing in this model because  $[H_0, H_1] = 0$ ), but merely due to the choice of the ensemble with the optimal number of doubly occupied sites. Finally we note that the axes in Fig. 1 of Ref. [1] should read u/4|t| and U/4|t|, respectively, to yield the correct Lieb-Wu result for the

Hubbard model in d=1. Even then the ground-state energy of the 1-d Hubbard model at U=0,  $E_{\text{Hubb}}^0/t = -4/\pi$ , does not coincide with the value from (5) for d=2 and U=0,  $E_g^0/t = -16/\pi^2$ .

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