Why one needs a functional RG to survive in a disordered world

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Abstract

In these proceedings, we discuss why Functional Renormalization is an essential tool to treat strongly disordered systems. More specifically we treat elastic manifolds in a disordered environment. These are governed by a disorder distribution, which after a finite renormalization becomes non-analytic, thus overcoming the predictions of the seemingly exact dimensional reduction. We discuss, how a renormalizable field theory can be constructed, even beyond 2-loop order. We then consider an elastic manifold imbedded in $N$ dimensions, and give the exact solution for $N \to \infty$. This is compared to predictions of the Gaussian replica variational ansatz, using replica symmetry breaking. Finally, the effective action at order $1/N$ is reported.

1 Introduction

In these proceedings we consider an elastic manifold in a random potential, as prototype for strongly disordered systems. Since for all these systems temperature is irrelevant, we will only treat zero temperature. The kind of systems we have in mind are domain walls in dirty magnets, contact lines, charge density waves, vortex lattices, to just mention a few. These results were obtained in collaboration with Pierre Le Doussal [1–11]. For lack of space we restrict our discussion to the equilibrium. Complementary material, especially for the depinning, can be found in the earlier review [12].

2 Physical realizations, model and observables

The simplest experimental realization is an Ising magnet. Imposing boundary conditions with all spins up at the upper and all spins down at the lower boundary (see figure 1), at low temperatures, a domain wall separates a region with spin up from a region with spin down. In a pure system at temperature $T = 0$, this domain wall is completely flat. Disorder can deform the domain wall, making it eventually rough again. Figure 1 shows, how the domain wall is described by a displacement field $u(x)$. Another example is the contact line of water (or liquid Helium), wetting a rough substrate. A realization with a 2-parameter displacement field $\vec{u}(\vec{x})$ is the deformation of a vortex lattice: the position of each vortex is deformed from $\vec{x}$ to $\vec{x} + \vec{u}(\vec{x})$. A 3-dimensional example are charge density waves.

All these models have in common, that they are described by a displacement field $x \in \mathbb{R}^d \rightarrow \bar{u}(x) \in \mathbb{R}^N$. For simplicity, we set $N = 1$, if not explicitly stated otherwise. After some initial coarse-graining, the energy $\mathcal{H} = \mathcal{H}_{\text{el}} + \mathcal{H}_{\text{DO}}$ consists out of two parts: the elastic energy $\mathcal{H}_{\text{el}}$ and the disorder energy $\mathcal{H}_{\text{DO}}$

$$\mathcal{H}_{\text{el}}[u] = \int d^d x \frac{1}{2} (\nabla u(x))^2, \quad \mathcal{H}_{\text{DO}}[u] = \int d^d x V(x, u(x))$$

We choose the disorder at the microscopic scale Gaussian, with correlations

$$V(u, x)V(u', x') := \delta^d(x - x') R(u - u').$$

The most interesting observable is the roughness-exponent $\zeta$, from the behavior of the correlation function

$$[u(x) - u(y)]^2 \sim |x - y|^{2\zeta}.$$  

Other observables are higher correlation functions or the free energy.
3 Dimensional reduction

There is a beautiful and rather mind-boggling theorem relating disordered systems to pure systems (i.e. without disorder), which applies to a large class of systems, e.g. random field systems and elastic manifolds in disorder. It is called dimensional reduction and reads as follows [14]:

**Theorem:** A $d$-dimensional disordered system at zero temperature is equivalent to all orders in perturbation theory to a pure system in $d - 2$ dimensions at finite temperature.

Let me give an example: The thermal expectation value for the 2-point function scales as $\langle |u(x) - u(y)|^2 \rangle \sim |x|^{2-d}$. Making the dimensional shift implied by dimensional reduction implies that the disorder-averaged 2-point function at zero temperature is

$$\langle (u(x) - u(0))^2 \rangle \sim x^{4-d} \equiv x^{2\zeta} \quad \text{i.e.} \quad \zeta = \frac{4-d}{2} .$$

(3.1)

We will see later that this is not true; but remains an important benchmark due to fact that the “theorem” is correct to all orders in the disorder strength and its moments (i.e. when expanding in $R''(0), R''''(0),$ a.s.o.).

4 The Larkin-length

To understand the failure of dimensional reduction, let us turn to an interesting argument given by Larkin [15]. He considers a piece of an elastic manifold of size $L$. If the disorder has correlation length $r$, and characteristic potential energy $\bar{\epsilon}$, this piece will typically see a potential energy of strength $E_{DO} = \bar{\epsilon}(Lr)^{d/2}$. On the other hand, there is an elastic energy, which scales like $E_{el} = cL^{d-2}$. These energies are balanced at the Larkin-length $L = L_c$ with $L_c = (\frac{\bar{\epsilon} r^d}{c})^{1/(4-d)}$. More important than this value is the observation that in all physically interesting dimensions $d < 4$, and at scales $L > L_c$, the membrane is pinned by disorder, whereas on small scales elastic energy dominates. This means that $d = 4$ is the upper critical dimension.

5 The functional renormalization group (FRG)

Let us now discuss a way out of the dilemma, posed by dimensional reduction: We would like to make an $\epsilon = 4 - d$ expansion. On the other hand, dimensional reduction tells us that the roughness is $\zeta = \frac{4-d}{2}$ (see (3.1)). Even though this is systematically wrong below four dimensions, it tells us correctly that at the critical dimension $d = 4$, where disorder is marginally relevant, the field $u$ is dimensionless. This means that having identified any relevant or marginal perturbation (as the disorder), we find
immediately another such perturbation by adding more powers of the field. We can thus not restrict ourselves to keeping solely the first moments of the disorder, but have to keep the whole disorder-distribution function $R(u)$. Thus we need a functional renormalization group treatment (FRG). Functional renormalization is an old idea going back to the seventies, and can e.g. be found in [16], by Wegner and Houghton. For disordered systems, it was first proposed in 1986 by D. Fisher [17]. Performing an infinitesimal renormalization, i.e. integrating over a momentum shell à la Wilson, leads to the flow of the first moments of the disorder, but have to keep the whole disorder-distribution function immediately another such perturbation by adding more powers of the field. We can thus not restrict ourselves to keeping solely the first moments of the disorder, but have to keep the whole disorder-distribution function $R(u)$. Thus we need a functional renormalization group treatment (FRG). 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want to integrate out a single degree of freedom, whose average position due to the elastic energy connecting it to its neighbors is $u$. This harmonic potential and the disorder term are represented by the parabola and the lowest curve on figure 3(a) respectively; their sum is the remaining curve. For a given disorder realization, the minimum of the potential as a function of $u$ is reported on figure 3(b). Note that it has non-analytic points, which mark the transition from one minimum to another. Taking the derivative of the potential leads to the force in figure 3(c). It is characterized by almost linear pieces, and shocks (i.e. jumps). Calculating the force-force correlator, the dominant contribution in its decay for small distances is due to the presence of shocks. Their contribution is proportional to their probability, itself proportional to the distance between the two observable points. This leads to $F(u)F(0) = F(0)^2 - c|u|$, with some numerical coefficient $c$.

## 7 Beyond 1 loop?

Functional renormalization has successfully been applied to a bunch of problems at 1-loop order. From a field theory, we however demand more. Namely that it allows for systematic corrections beyond 1-loop order; be renormalizable; and thus allows to make universal predictions. However, this has been a puzzle since 1986, and it has even been suggested that the theory is not renormalizable due to the appearance of terms of order $\epsilon^{3/2}$ [19]. Why is the next order so complicated? The reason is that it involves terms proportional to $R''(0)$. A look at figure 3 explains the puzzle. Shall we use the symmetry of $R(u)$ to conclude that $R''(0)$ is 0? Or shall we take the left-hand or right-hand derivatives, related by

$$R''(0^+) := \lim_{u \to 0^+} R''(u) = - \lim_{u \to 0^-} R''(u) =: -R''(0^-).$$

(7.1)

In the following, I will present the solution of this puzzle, at 2-loop order and large $N$. The latter approach allows for another independent control-parameter, and sheds further light on the cusp-formation.

## 8 Results at 2-loop order

For the flow-equation at 2-loop order, the result is [1,4,20,21]

$$\partial_\epsilon R(u) = (\epsilon - 4\zeta) R(u) + \zeta u R'(u) + \frac{1}{2} R''(u) R''(0) - R''(u) R''(0) + \frac{1}{2} (R'(u) - R''(0)) R''(u) R''(0) - \frac{1}{2} R''(0) R''(u).$$

(8.1)

The first line is the result at 1-loop order, already given in (5.1). The second line is new. The most interesting term is the last one, which involves $R''(0^+)^2$ and which we therefore call anomalous. The hard task is to fix the prefactor $(-\frac{1}{2})$. We have found five different prescriptions to calculate it: The slope-algorithm, recursive construction, reparametrization invariance, renormalizability, and potentiality [1,22]. For lack of space, we restrain our discussion to the last two ones. At 2-loop order the following diagram appears

leading to the anomalous term. The integral (not written here) contains a subdivergence, which is indicated by the box. Renormalizability demands that its leading divergence (which is of order $1/\epsilon^2$) be canceled by a 1-loop counter-term. The latter is unique thus fixing the prefactor of the anomalous term.

Another very physical demand is that the problem remain potential, i.e. that forces still derive from a potential. The force-force correlation function being $-R''(u)$, this means that the flow of $R''(0)$ has to be strictly 0. From (8) one can check that this does not remain true if one changes the prefactor of the last term in (8); thus fixing it.

Let us give some results for random-bond disorder (short-ranged potential-potential correlation function). For this, we have to solve (8.1) numerically, with the result $\zeta = 0.20829804 \epsilon + 0.00685862 \epsilon^2$. This compares well with numerical simulations, see figure 4.

| $d$ | one loop | two loop | estimate | simulation and exact |
|-----|----------|----------|-----------|----------------------|
| 3   | 0.208    | 0.215    | 0.215 ± 0.01 | 0.22 ± 0.01 [23] |
| 2   | 0.417    | 0.444    | 0.42 ± 0.02  | 0.41 ± 0.01 [23] |
| 1   | 0.625    | 0.687    | 0.67 ± 0.02  | 2/3                  |

Figure 4: Results for $\zeta$ in the random bond case.
calculated analytically as of this analytic solution: First of all, for long-range correlated disorder of the form w.r.t. This is a complicated nonlinear partial differential equation. It is therefore surprising, that one can find an analytic solution. (The trick is to write down the flow-equation for the inverse function of range correlated disorder, \(u\). We conclude that \(\int \partial \tilde{B}(x) \equiv -\frac{m\partial}{\partial m} \tilde{B}(x) = (\epsilon - 4\zeta)\tilde{B}(x) + 2\zeta x\tilde{B}'(x) + \frac{1}{2} \tilde{B}'(x)^2 - \tilde{B}'(x)\tilde{B}'(0)\) (10.2)

This is a complicated nonlinear partial differential equation. It is therefore surprising, that one can find an analytic solution. (The trick is to write down the flow-equation for the inverse function of \(\tilde{B}'(x)\), which is linear.) Let us only give the results of this analytic solution: First of all, for long-range correlated disorder of the form \(\tilde{B}'(x) \sim x^{-\gamma}\), the exponent \(\zeta\) can be calculated analytically as \(\zeta = \frac{1}{1+\gamma}\). It agrees with the replica-treatment in [25] and the 1-loop treatment in [19]. For short-range correlated disorder, \(\zeta = 0\). Second, it demonstrates that before the Larkin-length, \(\tilde{B}'(x)\) is analytic and thus dimensional reduction holds. Beyond the Larkin length, \(\tilde{B}''(0) = \infty\), a cusp appears and dimensional reduction is incorrect. This shows again that the cusp is not an artifact of the perturbative expansion, but an important property even of the exact solution of the problem (here in the limit of large \(N\)).

9 Finite \(N\)

Up to now, we have studied the functional RG in two cases: For one component \(N = 1\) and in the limit of a large number of components, \(N \to \infty\). The general case of finite \(N\) is more difficult to handle, since derivatives of the renormalized disorder now depend on the direction, in which this derivative are taken. Define amplitude \(u := |\tilde{u}|\) and direction \(\hat{u} := \tilde{u}/|\tilde{u}|\) of the field. Then deriving the latter variable leads to terms proportional to \(1/u\), which are diverging in the limit of \(u \to 0\). This poses additional problems in the calculation, beyond the case \(N = 1\). At 1-loop order everything is well-defined [19]. We have found a consistent RG-equation at 2-loop order (see [12] and unpublished):

The fixed point equation has to be integrated numerically, order by order in \(\epsilon\). The result, specialized to directed polymers, i.e. \(\epsilon = 3\) is plotted on figure 5. We see that the 2-loop corrections are rather big at large \(N\), so some doubt on the applicability of the latter down to \(\epsilon = 3\) is advised. However both 1- and 2-loop results reproduce well the two known points on the curve: \(\zeta = 2/3\) for \(N = 1\) and \(\zeta = 0\) for \(N = \infty\). The latter result has been given in section 10 Via the equivalence [24] of the directed polymer problem in \(N\) dimensions treated here and the KPZ-equation of non-linear surface growth in \(N\) dimensions, we conclude that \(d \approx 2.4\) is the upper critical dimension of KPZ.

10 Large \(N\)

In the last section, we have discussed renormalization in a loop expansion, i.e. expansion in \(\epsilon\). In order to independently check consistency it is good to have a non-perturbative approach. This is achieved by the large-\(N\) component field \(\tilde{u}\). We then calculate the free energy in presence of a source \(j\), and finally the effective action \(\Gamma(\tilde{u})\). This poses some doubt on the applicability of the latter down to \(\epsilon = 3\) is advised. However both 1- and 2-loop results reproduce well the two known points on the curve: \(\zeta = 2/3\) for \(N = 1\) and \(\zeta = 0\) for \(N = \infty\). The latter result has been given in section 10 Via the equivalence [24] of the directed polymer problem in \(N\) dimensions treated here and the KPZ-equation of non-linear surface growth in \(N\) dimensions, we conclude that \(d \approx 2.4\) is the upper critical dimension of KPZ.
11 Relation to Replica Symmetry Breaking (RSB)

There is another treatment of the limit of large $N$ given by Mézard and Parisi [25]. They make a Gaussian variational ansatz of the form

$$H_{\mathbf{g}}[\mathbf{u}] = \frac{1}{2T} \sum_{a=1}^{n} \int_{x} \bar{u}_{a}(x) (-\nabla^{2} + m^{2}) \bar{u}_{a}(x) - \frac{1}{2T^{2}} \sum_{a,b=1}^{n} \sigma_{ab} \bar{u}_{a}(x) \bar{u}_{b}(x),$$

(11.1)

which becomes exact for $N \to \infty$. The art is to make an appropriate ansatz for $\sigma_{ab}$. The simplest possibility, $\sigma_{ab} = \sigma$ for all $a \neq b$ reproduces the dimensional reduction result, which breaks down at the Larkin length. Beyond that scale, a replica-symmetry broken (RSB) ansatz for $\sigma_{ab}$ is necessary, of the form $\sigma_{ab} = \sigma(0)$. Parisi has shown that this infinitely often replica-symmetry broken matrix can be parameterized by a function $\sigma(z)$ with $z \in [0,1]$ where $z = 0$ describes distant states, whereas $z = 1$ describes nearby states. The solution of the large-$N$ saddle-point equations leads to the curve depicted in figure 11 Knowing it, the 2-point function is given by

$$\langle u_{k} u_{-k} \rangle = \left. \frac{1}{k^{2} + m^{2}} \right|_{\chi_{ab}} + \int_{0}^{1} \frac{dz}{k^{4} + [\sigma(z) + m^{2}]}.$$

(11.2)

What is the relation between the two approaches, which both pretend to calculate the same 2-point function? Comparing the analytical solutions, we find that the 2-point function given by FRG is the same as that of RSB, if in the latter expression we only take into account the contribution from the most distant states, i.e. those for $z$ between $0$ and $z_{m}$ (see figure 6). To understand why this is so, we have to remember that the two calculations were done under quite different assumptions: In contrast to the RSB-calculation, the FRG-approach calculated the partition function in presence of an external field $j$, which was then used to give via a Legendre transformation the effective action. Even if the field $j$ is finally turned to 0, the system will remember its preparation, as is the case for a magnet.

By explicitly breaking the replica-symmetry through an applied field, all replicas will settle in distant states, and the close states from the Parisi-function $[\sigma(z) + m^{2}]$ (which describes spontaneous RSB) will not contribute. However, we found that the full RSB-result can be reconstructed by remarking that the part of the curve between $z_{m}$ and $z_{c}$ is independent of the infrared cutoff $m$, and then integrating over $m$ [2] ($m_{c}$ is the mass corresponding to $z_{c}$):

$$\langle u_{k} u_{-k} \rangle_{RSB}^{IR−cutoff} = \frac{\tilde{R}_{m}(0)}{m^{2}} + \int_{m}^{m_{c}} \frac{d \tilde{R}_{m}(0)}{\mu^{2}} + \frac{1}{m_{c}^{2}} = \frac{1}{m^{2}}.$$

(11.2)

We also note that a similar effective action has been proposed in [18]. While it agrees qualitatively, it does not reproduce the correct FRG 2-point function, as it should.

12 Corrections at order $1/N$

In a graphical notation, we find [11]

$$\delta B^{(1)} = \text{graphical expression} + \text{other terms},$$

+ $T \left( \text{graphical expression} + \text{other terms} \right)$

+ $T^{2} \left( \text{graphical expression} + \text{other terms} \right)$

(12.1)

$$\boldsymbol{\sigma} = B'(\chi_{ab}) (1 - 4A_{d} I_{z}(p) B''(\chi_{ab}))^{-1}, \quad \chi_{ab} = B(\chi_{ab}),$$

(12.2)
where the explicit expressions are given in [11]. By varying the IR-regulator, one can derive a $\beta$-function at order $1/N$, see [11]. At $T = 0$, it is UV-convegent, and should allow to find a fixed point. We have been able to do this at order $\epsilon$, showing consistency with the 1-loop result, see section 9. Other dimensions are more complicated.

A $\beta$-function can also be defined at finite $T$. However since temperature is an irrelevant variable, it makes the theory non-renormalizable, i.e. in order to define it, one must keep an explicit infrared cutoff. These problems will be treated in a forthcoming publication.

13 Perspectives

Other interesting problems have been treated by the above methods, especially dynamic problems (see [12] for a review); and many more are now in reach. Some open points have already been raised in these notes, others are the strong disorder phase of random field problems, or whether FRG can also be applied to spin-glasses. We have to leave these problems for future research and as a challenge for the reader to plunge deeper into the mysteries of functional renormalization.

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