Flow equation holography

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The Ryu-Takayanagi conjecture establishes a remarkable connection between quantum systems and geometry. Specifically, it relates the entanglement entropy to minimal surfaces within the setting of AdS/CFT correspondence. This Letter shows how this idea can be generalised to generic quantum many-body systems within a perturbative expansion where the region whose entanglement properties one is interested in is weakly coupled to the rest of the system. A simple expression is derived that relates a unitary disentangling flow in an emergent RG-like direction to the min-entropy of the region under consideration. Explicit calculations for critical free fermions in one and two dimensions illustrate this relation.

AdS/CFT correspondence [1,2] as an explicit realization of the holographic principle [3,4] has become a powerful tool for understanding properties of strongly interacting quantum field theories. It not only establishes a connection between certain classes of gauge theories in the ’t Hooft limit and classical theories of gravity with an additional emergent RG-like dimension, but also led to a new way of thinking about quantum field theories. For example, the Ryu-Takayanagi conjecture [5–7] demonstrates a remarkable relation between the entanglement entropy of a region \( A \subset \mathbb{R}^d \) in a \( d+1 \) dimensional conformal field theory and the static minimal surface in the dual \( d+2 \) dimensional AdS-space with a boundary given by \( \partial A \). This is important both conceptually and also from a practical point of view since the calculation of entanglement entropies is challenging even for ground states of simple systems, with few analytical or numerical tools available, especially for \( d > 1 \). It has become increasingly clear in the past decade that the entanglement entropy plays a fundamental role in fields as diverse as quantum information theory [8], the efficient simulation of quantum systems [9,10] or topological phases of matter [11].

Therefore the insight described by the Ryu-Takayanagi conjecture spurred a lot of activity to find similar relations between entanglement properties in \( d+1 \) dimensions and geometry in \( d+2 \) dimensions beyond the limitation to AdS/CFT-theories. Especially the geometric picture underlying the multi-scale entanglement renormalization ansatz (MERA) [12] and its continuous version cMERA [13] bears strong resemblance to AdS/CFT-correspondence and the Ryu-Takayanagi conjecture [14], while in principle being applicable for a wide class of quantum many-body systems. Still it remains to be explored to what extent these tools can also be used for a quantitative calculation of entanglement properties.

Motivated by these ideas this Letter develops a systematic procedure which connects the entanglement properties of eigenstates of a generic quantum many-body Hamiltonian to a disentangling flow in an emergent RG-like dimension. This is done by working in a controlled limit where the subregion \( B \) whose entanglement we are interested in is weakly coupled to the rest of the system, see Fig.1 I denote the relative strength of this coupling by \( g \) with \( |g| \ll 1 \) being the weak-link limit. \( g = 1 \) will correspond to translation-invariant systems in the models discussed explicitly later on.

For a bipartite Hilbert space \( \mathcal{H} = \mathcal{H}_A \otimes \mathcal{H}_B \) the entanglement of region \( B \) for a pure state \( |\psi\rangle \in \mathcal{H} \) is described by its Rényi entropies

\[
S^{(q)} = \frac{1}{1-q} \ln \text{tr}_{\mathcal{H}_B} \rho_B^q ,
\]

where \( q \) is a positive real number and \( \rho_B = \text{tr}_{\mathcal{H}_A} |\psi\rangle \langle \psi| \) is the reduced density matrix of subsystem \( B \). The limit \( q \to 1 \) yields the von Neumann entanglement entropy. The challenge for the analytic calculation of entanglement entropies of gapless models (and the reason why relatively little is known analytically even in one dimension apart from free theories or conformal field theories) is that the eigenvalues of the reduced density matrix behave non-perturbatively with the size of region \( B \). This is even true in the weak-link limit and one also finds non-perturbative behavior in \( g \) [15].

As worked out below, this problem can be resolved for the Rényi entropy in the limit \( q \to \infty \), which is also called
the min-entropy or single-copy entropy [10]. The min-entropy is only determined by the largest eigenvalue $p_{\text{max}}$ of the reduced density matrix $\rho_B$, 

$$S^{(\text{min})} = \lim_{q \to \infty} S^{(q)} = -\ln p_{\text{max}} .$$

(2)

We will see that $p_{\text{max}}$ has a specific product structure emerging from the disentangling flow that leads to a non-perturbative result for $p_{\text{max}}$ even if the individual factors are only known perturbatively in $g$. The goal of this Letter is to give an explicit expression for $S^{(\text{min})}$ which is correct in order $g^2$ based on a disentangling flow without making strong AdS/CFT-like restrictions on the class of permissible systems. The focus on $S^{(\text{min})}$ as opposed to the von Neumann entropy is also one main difference to prior work on disentangling flows [16], which allows me to give an analytic justification in the weak-link limit. It should also be mentioned that the min-entropy has an important quantum information theory interpretation, namely as the amount of distillable entanglement from a single copy of the quantum state $|\psi\rangle$, as opposed to the von Neumann entropy which requires many identically prepared quantum states [10].

The key step in my approach is the construction of an RG-like disentangling flow described by unitary transformations $U(u)$. In this Letter we are interested in the entanglement properties of eigenstates of Hamiltonians, $H|\psi\rangle = E|\psi\rangle$. In order to find a disentangling flow for $|\psi\rangle$ we unitarily transform the Hamiltonian $H(u) \overset{\text{def}}{=} U(u) H U(d) u$ in such a way that $H(u = \infty)$ only contains terms which act locally in $\mathcal{H}_A$ or $\mathcal{H}_B$, and no longer contains terms which generate entanglement across the links between regions $A$ and $B$. Since the unitarily transformed state $|\psi(u)\rangle = U(u)|\psi\rangle$ is always an eigenstate of $H(u)$, this implies that $|\psi(u = \infty)\rangle$ must be a disentangled state since it is an eigenstate of $H(u = \infty)$.

A unitary flow can be obtained following the flow equation method [17, 18] which generates $U(u)$ from a sequence of infinitesimal unitary transformations with an antihermitean generator $\eta(u)$

$$\frac{dH(u)}{du} = [\eta(u), H(u)]$$

(3)

and initial condition $H(u = 0) = H$. The unitary transformation is obtained as an $u$-ordered exponential

$$U(u) = T_u \exp \left( \int_0^u du' \eta(u') \right)$$

(4)

where the larger values of $u'$ are ordered to the left. A flow with the desired properties is obtained by splitting up the Hamiltonian as

$$H(u) = H_A(u) + H_B(u) + H_{\text{ent}}(u)$$

(5)

where only $H_{\text{ent}}(u)$ contains terms that create entanglement by linking regions $A$ and $B$. The generator is defined as

$$\eta(u) \overset{\text{def}}{=} 2\Lambda^2 e^{2u} [H(u), H_{\text{ent}}(u)]$$

(6)

where $\Lambda$ is the ultraviolet cutoff. One can show that under very general conditions this choice of the generator leads to a contracting flow, $\frac{d}{du} \text{tr}(H^2_{\text{ent}}(u)) \leq 0$ and $\eta(u = \infty) = 0$ [17, 18]. While this does not guarantee that $H_{\text{ent}}(u = \infty)$ vanishes, it commutes with $H_A$ and $H_B$ and does not contribute to entanglement between regions $A$ and $B$ anymore (in the sense of not affecting the leading behavior of the entanglement entropy with the size of $B$). Previous applications of the flow equation method were such that an interaction part in the Hamiltonian was eliminated by a sequence of infinitesimal unitary transformations (see e.g. Refs. [17, 20] and further references therein). The main difference in this Letter is that I use flow equations to disentangle a Hamiltonian, that is to eliminate couplings between two regions for flow parameter $u \to \infty$. In practise Eqs. (3) and (6) lead to coupled systems of ordinary differential equations, which for interacting many-body systems generically need to be truncated in order to render the hierarchy finite. This can be done easily in the weak-link limit since by definition $H_{\text{ent}}$ is proportional to $g$, therefore the flow equations (3) and (6) can be organised in powers of $g$. The resulting systems of differential equations can be solved either numerically or (usually approximately) analytically (more below).

Notice that $u$ can be thought of as a logarithmic energy scale, $u = \ln(A/E)$, and that the generator $\eta(u)$ is dimensionless [21]. Matrix elements in $H_{\text{ent}}$ that couple states with an energy difference $E$ are unitarily transformed away at flow scale $u$. So large energy differences are eliminated first (for small $u$) and smaller energy differences later along the flow, which makes the flow equation method RG-like. Energy differences correspond to length scales and for translation-invariant systems (meaning: translation-invariant apart from the weak links) the unitary transformation acts on sites that are further and further apart from the boundary between regions $A$ and $B$ as $u$ increases. This means that the Hilbert spaces and the unitary transformation can be factorized

$$\mathcal{H}_{A,B} = \bigotimes_i \mathcal{H}^{(i)}_{A,B} , \quad U(u) = \prod_i U^{(i)}(u)$$

(7)

such that $U^{(i)}$ only acts nontrivially on $\mathcal{H}^{(i)}_{A,B} \otimes \mathcal{H}^{(i)}_{B,B}$. The index $i$ does not need to refer to single sites, $\mathcal{H}^{(i)}_{A,B}$ can contain many sites as long as a perturbative argument for $U^{(i)}$ remains valid (see below).

If one reconstructs $|\psi\rangle$ from the disentangled state $|a_0\rangle \otimes |b_0\rangle$, one therefore has a structure as depicted in Fig. 2. Each individual $U^{(i)}(\infty)$ is close to the identity in the weak-link limit, hence the component $|a_0\rangle \otimes |b_0\rangle$ of the total wavefunction gets multiplied by a factor $\lambda^{(i)}_{\text{max}} = 1 - O(g^2)$ under the action of $U^{(i)}(\infty)$. In addition, $U^{(i)}(\infty)$ generates a component orthogonal to $|a_0\rangle \otimes |b_0\rangle$, which however cannot contribute to the coef-
ficient in front of \(|a_0\rangle \otimes |b_0\rangle\) anymore in later steps. This implies that the Schmitt decomposition of \(|\psi\rangle\) takes the form \(|\psi\rangle = \sum_j \lambda_j (a_j \otimes b_j)\) with the largest Schmitt coefficient \(\lambda_0 = \prod_i \lambda^{(i)}_{\text{max}}\). We will show \(\lambda^{(i)}_{\text{max}} = 1 - \alpha^{(i)} g^2\) with coefficients \(\alpha^{(i)}\) that can be calculated perturbatively, therefore

\[
S^{(\text{min})} = -\ln p_{\text{max}} = -\ln \lambda_0^2 = -2 \ln \prod_i \lambda^{(i)}_{\text{max}}
\]

\[
= -2 \sum_i \ln \lambda^{(i)}_{\text{max}} = 2g^2 \sum_i \alpha^{(i)} + O(g^4) . \quad (8)
\]

In this manner one avoids calculating \(p_{\text{max}}\) with its non-perturbative behavior by making use of the product structure for the largest Schmitt coefficient depicted in Fig. 2. It is sufficient to find the individual \(\lambda^{(i)}_{\text{max}}\) perturbatively in the weak-link limit. This is a straightforward exercise in second order perturbation theory: If one weakly perturbs a Hamiltonian \(H_0, H_g = H_0 + gH_{\text{int}}\), then the overlap of the perturbed eigenstate with the unperturbed eigenstate can be written as

\[
|\psi_g \rangle |\psi_0\rangle = 1 - \int_0^\infty du \left( - \langle \psi_0 | \eta^2(u) | \psi_0 \rangle \right) + O(g^3) \quad (9)
\]

with the flow generated from (3) and \(\eta(u) \equiv 2\Lambda^{-2} e^{2\alpha \delta t} [H_g(u), gH_{\text{int}}(u)]\) like in (6). This implies for (8)

\[
S^{(\text{min})} = -2 \sum_i \int_0^\infty du \left( \langle \psi | \eta^{(i)}(u) \rangle^2 | \psi \rangle \right) + O(g^3) \quad (10)
\]

where \(\eta^{(i)}\) is the infinitesimal generator corresponding to \(U^{(i)}\). Since by construction \(\langle \psi | \eta^{(i)}(u) \eta^{(j)}(u) | \psi \rangle \propto \delta_{ij}\), we may define

\[
S^{(\text{feq})} \overset{\text{def}}{=} -2 \int_0^\infty du \langle \psi | \eta^2(u) | \psi \rangle \quad (11)
\]

and have shown

\[
S^{(\text{min})} = S^{(\text{feq})} + O(g^3) . \quad (12)
\]

These two formulas constitute the main result of this Letter. They connect the min-entropy to the integral over the flow equation disentangling density measured by \(D(u) \overset{\text{def}}{=} -\langle \psi | \eta^2(u) | \psi \rangle\) on a logarithmic energy scale \((D(u) \geq 0\) since \(\eta(u)\) is antithermitean).

In order to demonstrate this approach we study critical free fermions in \(d = 1\) and \(d = 2\) explicitly. This will permit us to compare with various exact analytical or well-controlled numerical results. In \(d = 1\) we consider the Hamiltonian

\[
H = - \sum_{i=-\infty}^{\infty} t_i \left( c_i^\dagger c_{i+1} + \text{h.c.} \right) \quad (13)
\]

with \(t_0 = t_{\ell} = g\) and \(\forall i \neq 0, \ell\quad t_i = 1\). We are interested in the ground state entanglement entropy of the lattice sites \(\{1, \ldots, \ell\}\) with the rest of the chain. For the translation-invariant case \(g = 1\) the behavior is logarithmic in \(\ell\) with a universal prefactor [22–24] independent of the filling, \(S(q) = \frac{1-q}{6} \ln \ell + O(\ell^0)\). Our goal is now to extract this prefactor in the min-entropy as a function of the link strength \(g\)

\[
S^{(\text{min})} = \kappa(g) \ln \ell + O(\ell^0) \quad (14)
\]

using the disentangling flow. We do this by considering [13] with one weak link \((t_0 = g\) and \(\forall i \neq 0, \ell\quad t_i = 1\) and following the disentangling flow separating the left half line from the right half line. Initially \(H_{\text{ent}}(u = 0) = -g(c_j^\dagger c_1 + \text{h.c.})\) and during the flow \(H_{\text{ent}}(u)\) in (5) is defined to consist of the hopping terms in \(H(u)\) connecting the sites \(\{-\infty, \ldots, 0\}\) to the sites \(\{1, \ldots, \infty\}\). The resulting system of differential equations (3) is solved numerically (without approximations or truncations). After an
integrate up to values $u$ gives $\kappa$ and does not affect the logarithmic behavior. The initial transient one finds that the flow equation disentangling density $D(u)$ becomes constant. An upper cut-off in (11) is set by the observation that $H_{\text{ent}}(u)$ becomes increasingly non-local during the flow: The initial nearest neighbor hopping in $H_{\text{ent}}(u = 0)$ turns into hopping terms between sites with separation $d$ (compare Fig. 2) at the scale $u = \ln d$. Since we are interested in decoupling an interval of length $\ell$ in (13) this implies we should only integrate up to values $u^* = \ln \ell$ in (11). Taking into account that the entanglement contributions from the two boundaries in (13) add up (for sufficiently large $\ell$) this gives $\kappa(g) = 4D$. Notice that a proportionality factor $p$ in $u^* = \ln(p \ell)$ only contributes to the $O(\ell^0)$-term in (14) and does not affect the logarithmic behavior.

The results are shown in Fig. 3 as a function of the link strength $g$ and for various fillings $\kappa(g)/g^2$ is depicted since $\kappa(g) \propto g^2$ for small $g$. As expected from (12) the flow equation results agree with exact diagonalization results for the min-entropy for small $g$ (23). Not surprisingly there are differences between $S^{\text{min}}$ and $S^{(\text{ten})}$ for larger values of $g$. For the translation-invariant case $g = 1$ the exact diagonalization results become independent from the filling as known from universality with $\kappa = 1/6$. Interestingly, within numerical accuracy the logarithmic prefactor in the flow equation expression $S^{(\text{ten})}$ also becomes independent from filling at this point.

Next we test the flow equation disentangling approach for a two-dimensional system, namely for the ground state of free fermions with nearest neighbor hopping ($t = 1$) on a square lattice which is infinite in the $x$-direction and has periodic boundary conditions in the $y$-direction (circumference $N_x \gg 1$ sites). We are interested in the min-entropy of a stripe with dimensions $\ell \times N_y$, which is coupled to the rest of the system with hopping matrix elements $g$. The general behavior is an area law with logarithmic corrections (10), that is $\kappa(g) \propto N_y$ in (14). This comes out naturally from (11) since $-\langle \psi | \eta^2(u) | \psi \rangle \propto N_y$ for large $N_y$. Results from the numerical solution of the flow equations are depicted in Fig. 4 and compared with exact diagonalization (26). For small $g$ one again finds perfect agreement as expected from (12).

In the one-dimensional case one can also interpret (11) from the point of view of minimal curves like in the Ryu-Takayanagi conjecture (5). With the emergent dimension $z = \Lambda^{-1} e^u$ one can verify that the leading behavior of (11) can also be written as the length of the minimal curve, $S^{(\text{ten})} = \min_{\mathcal{C}} L(\mathcal{C})$, for given UV-boundary conditions $\partial \mathcal{C} = \partial B$ with respect to the metric
ds = 4D (\ln(z\Lambda))^2 dz^2 + dx^2 \quad (15)

For the critical fermions discussed before this is just an AdS$_2$-metric since the flow equation disentangling density is constant on a logarithmic scale, $D(u) = \text{const}$. The spatial part $z^{-2} dx^2$ in (15) effectively describes the increasing non-locality of $H_{\text{ent}}(u)$ during the flow and thereby introduces the IR-cutoff $u^*$ once the length scale generated by the disentangling flow exceeds the size of the interval $B$ (here assuming that the total system is infinite). The minimal curve condition in the Ryu-Takayanagi conjecture finds a natural explanation in the flow equation framework since IR-cutoffs in the disentangling flow at a given boundary point are naturally set by the next boundary point closest to it. Consider for example a tight-binding chain (13) consisting of $N_x \gg 1$ lattice sites with periodic boundary conditions split up in two intervals with $\ell_A$ and $\ell_B$ sites, $\ell_A + \ell_B = N_x$. Then (11) gives

$$S^{(\text{ten})} = \kappa(g) \ln \min(\ell_A, \ell_B) + O(\ell^0) \quad (16)$$

which agrees with the minimal curve prescription using (15) and the known exact behavior (28).

In summary, I have introduced a method for calculating entanglement entropies for eigenstates of generic many-body Hamiltonians from a disentangling flow in an emergent RG-like direction (11). In the weak-link limit we have seen (12) that this procedure gives the min-entropy. While the specific examples discussed in this Letter were ground states of free fermion systems, the method also works for excited states and interacting systems. Notice that results in the weak-link limit can also be obtained analytically, although I have consistently used numerical methods for the solution of (3) in order to show results for all link strengths.
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[25] At half filling there is an exact analytical result for arbitrary $g$ [15], which is plotted as the dashed line in Fig. 3. It serves as a test for my exact diagonalization routine; its results agree within numerical precision as can be seen in Fig. 3.
[26] For the translation-invariant case $g = 1$ there is an analytic result [27] which has been used to validate my exact diagonalization numerics for $d = 2$. The agreement is excellent for all $\mu$ in Fig. 4 such that the differences are not visible on the scale of the plot.
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