

LETTER TO THE EDITOR

Smooth adiabatic evolutions with leaky power tails

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Abstract. Adiabatic evolutions *with* a gap condition have, under a range of circumstances, exponentially small tails that describe the leaking out of the spectral subspace. In general, adiabatic evolutions *without* a gap condition do not seem to have this feature. This is a known fact for eigenvalue crossing. We show that this is also the case for eigenvalues at the threshold of the continuous spectrum by considering the Friedrichs model.

1. Introduction

Adiabatic theorems describe the solutions of initial-value problems where the Hamiltonian generating the evolution depends slowly on time. In quantum mechanics the description is in terms of spectral information of the instantaneous Hamiltonian. A few basic references on various types of adiabatic theorems are [4, 7, 10, 12, 15, 16].

To formulate the problem more precisely it is convenient to replace the physical time t by the scaled time $s = t/\tau$. One is then concerned with the solution of the initial value problem

$$i\dot{\psi}_\tau(s) = \tau H(s)\psi_\tau(s) \quad (1)$$

with initial data $\psi_\tau(0) \in \text{Range } P(0)$. $P(s)$ is a spectral projection for $H(s)$, a self-adjoint Hamiltonian which depends sufficiently smoothly on s . ψ_τ is a vector-valued function and the adiabatic limit is the limit of large τ . Suppose, for the sake of concreteness, that $\dot{H}(s)$ is compactly supported on $[0, 1]$. As s runs on the interval $[0, 1]$, $H(s)$ evolves slowly in *physical* time for a *long* interval of time. The total variation of $H(s)$ is finite and not necessarily small.

Adiabatic theorems fall into two baskets: those that describe the solutions for *all* times, including times $s \in [0, 1]$, and those that characterize the solutions at large times $s > 1$, where the Hamiltonian is time independent again. Interestingly, they give more precision for long times. We call the first basket, the one that applies to all times, *uniform*, the second is called the *long-time* basket.

Adiabatic theorems can also be put in two other baskets: those that satisfy gap conditions and those that do not. By a gap condition we mean a spectral condition on the Hamiltonian $H(s)$ so that a finite gap separates the spectral subspace of the initial data ($= \text{Range } P(s)$), from the rest of the spectrum for all times s .

A representative result from the uniform basket is [1,4] presented in the following theorem.

Theorem 1. Suppose that $P(s)$ is smooth finite-rank spectral projection, for a bounded, smooth Hamiltonian $H(s)$. Then, the evolution of the initial state $\psi_\tau(0) \in \text{Range } P(0)$, is such that

$$\text{dist}(\psi_\tau(s), \text{Range } P(s)) \leq o(1) \tag{2}$$

for all $s \geq 0$. If, in addition, $P(s)$ satisfies a gap condition then

$$\text{dist}(\psi_\tau(s), \text{Range } P(s)) \leq O(1/\tau) \tag{3}$$

for all $s \geq 0$.

Remark 1. By $o(1)$ we mean a term that vanishes as $\tau \rightarrow \infty$.

Remark 2. Schrödinger operators constitute a class of application not covered by the theorem because of the assumption that $H(s)$ is bounded. This restriction can be lifted by standard machinery, see e.g. [1, 4]. We chose not to enter into this because the essence of the adiabatic evolution is an infrared problem which is largely divorced from the issue of unboundedness of the generator (an ultraviolet problem).

A characteristic result which lies both in the long-time basket and in the gap condition basket is [3, 6, 9, 11, 13, 14, 17] presented in the following theorem.

Theorem 2. Suppose that $H(s)$ is a smooth, bounded, and self-adjoint with $\dot{H}(s)$ supported on $[0, 1]$. And suppose that $P(s)$ is separated by a finite gap from the rest of the spectrum. Then, the evolution of the initial state $\psi_\tau(0) \in \text{Range } P(0)$, is such that for any $s > 1$,

$$\text{dist}(\psi_\tau(s), \text{Range } P(s)) = o\left(\frac{1}{\tau^n}\right) \tag{4}$$

for all $n \geq 0$.

Remarks.

- (i) There is, in general, no uniformity in n ; the term on the right-hand side is of order $\frac{c_n}{\tau^n}$, where c_n are allowed to grow rapidly with n .
- (ii) In the case where $\dot{H}(s)$ is an analytic family which decays at infinity, the leaking at $s = \infty$ is exponentially small.
- (iii) In the case of a gap, the distinction between the uniform and the long-time basket has an analogue in integrals. Suppose that $g(s)$ is real, smooth, strictly monotonic and compactly supported in $[0, 1]$. Then

$$\int_{-\infty}^s g(t)e^{it\tau} dt = \begin{cases} o(1/\tau^n) & \text{if } s > 1 \text{ for all } n \\ O(1/\tau) & \text{if } s \in [0, 1]. \end{cases}$$

The gapless basket seems to have no analogue in integrals, as we shall see.

What can one say about adiabatic theorems that lie both in the long-time and gapless baskets? In the uniform adiabatic theorem 1, the price one pays for the absence of a gap is, in general, loss of information on the rate at which the adiabatic limit is approached. The question we address here is what is the price one pays in the long-time adiabatic theorem, for the absence of a gap.

Adiabatic theorems without a gap condition are normally associated with two distinct settings: those that describe eigenvalue crossings and those associated with an eigenvalue at the threshold of the essential spectrum.

Hagedorn [8] studied adiabatic theorems for crossing eigenvalues. He showed that *there is no improvement in the adiabatic theorem in the long-time limit*. In particular, the leaking is always of order $\frac{1}{\sqrt{\tau}}$ for linear crossing.

Here we consider the cases where the absence of a gap is associated with an eigenvalue at the threshold of the continuous spectrum. The results of Hagedorn for crossing eigenvalues do not shed light on this case[†].

To investigate this we consider the Friedrichs model [5] with an eigenvalue at the threshold of the continuous spectrum. The model is parametrized by a real parameter $\beta > 0$ related to the behaviour of the density of states at low energies. Our main result is[‡]

$$\text{dist}(\psi_\tau(s), \text{Range } P(s)) = \begin{cases} O(\tau^{-\beta}) & 1 < \beta < 2 \quad \text{and} \quad s > 1 \\ O(\tau^{-1}) & 1 < \beta < 2 \quad \text{and} \quad s \in [0, 1]. \end{cases} \quad (5)$$

Unlike the gap basket, the long-time behaviour in the Friedrichs model indeed has power tails, and unlike the crossing basket, there is an improvement in the rate of decay at long times.

The main application of adiabatic theorems with eigenvalues at threshold is to the theory of atoms in radiation field, where the photon field eliminates spectral gaps. The absence of a gap makes the adiabatic theory of atoms in radiation field qualitatively different from quantum mechanics. Two distinct processes lead to the error in adiabatic evolutions. The first is photon production and the second is atomic excitation. Folk wisdom holds that the leaking due to atomic excitations has exponential tails. A simple argument suggests that the leaking due to photon production is of order $1/\tau$. The argument goes as follows: the power radiated by a dipole is, by classical electrodynamics,

$$\frac{2}{3} \frac{(d\dot{)}^2}{c^3}$$

where d is the dipole moment and c the velocity of light. It follows that the number of radiated photons (assuming characteristic frequency of order $O(1/\tau)$), is of order

$$\left(\frac{\alpha^{3/2}}{\tau}\right)^2$$

in atomic units[§].

2. Adiabatic evolutions

A basic strategy for studying adiabatic evolutions, introduced by Kato [12], is to compare the true evolution, $U_\tau(s)$ with a fictitious one, $U_a(s)$, which respects the spectral splitting. That is

$$P(s) = U_a(s) P U_a^\dagger(s). \quad (6)$$

We shall denote, by a subscript, the generator and the evolution of such a fictitious dynamics, i.e.

$$i\dot{U}_a(s) = \tau H_a(s) U_a(s) \quad U_a(0) = 1 \quad (7)$$

while for the true dynamics

$$i\dot{U}_\tau(s) = \tau H(s) U_\tau(s) \quad U_\tau(0) = 1. \quad (8)$$

[†] At crossing the spectral projection is discontinuous while for the eigenvalue at threshold that we consider the spectral projection is smooth for all times.

[‡] All estimates should be considered in an asymptotic sense with respect to $\tau \rightarrow \infty$.

[§] Recall that in atomic units $\alpha = 1/c$.

It was shown by Kato that H_a satisfies the commutator equation

$$\tau[H_a(s), P(s)] = i[\dot{P}(s), P(s)]. \quad (9)$$

This commutator equation does not have a unique solution, and different choices of H_a can be made. Kato chose

$$H_K = \frac{i}{\tau}[\dot{P}(s), P(s)]. \quad (10)$$

This generates a geometric evolution so that, in particular, $U_K(s)$ is independent of τ . It is a convenient choice for proving a uniform adiabatic theorem when $P(s)$ is the kernel of $H(s)$. It is inappropriate for generating a systematic adiabatic expansion.

A more effective choice, introduced in [2], is

$$H_{AD} = H(s) + \frac{i}{\tau}[\dot{P}(s), P(s)]. \quad (11)$$

This generator is close to the Hamiltonian and it can be used to prove stronger adiabatic theorems than those handled by Kato, and it can also be used to generate an adiabatic expansion and prove long-time adiabatic theorems [13]. However, it is often difficult to use this evolution for concrete computations because there is no explicit formula for $U_{AD}(s; \tau)$. This is the course one should follow when one wants to show fat leaky tails where one needs a *lower bound* on the part that leaks. For this reason we require yet another choice of an adiabatic evolution that more readily lends itself to explicit computation.

In the special case when the time dependence enters through the unitary family

$$H(s) = V(s)HV^\dagger(s) \quad (12)$$

the evolution in the rotating frame is both adiabatic and explicit:

$$U_r(s) = V(s) \exp(-i\tau s H). \quad (13)$$

It is generated by

$$H_r(s) = H(s) + \frac{i}{\tau}\dot{V}(s)V^\dagger(s) \quad (14)$$

and like H_{AD} is also close to $H(s)$. This evolution turns out to be inferior to H_{AD} when the game is to prove *upper bounds* on the leaky tails, but, it is a useful evolution in estimating fat leaky tails where the game is to obtain lower bounds. The best of all worlds occurs when H_{AD} and H_r coincide. This will turn out to be the case in the Friedrichs model we consider and this is what makes the analysis of this case simple.

3. The wave operator

To compare the true dynamics with the fictitious dynamics one introduces the wave operator

$$\Omega_{AD}(s; \tau) = U_{AD}^\dagger(s; \tau)U_\tau(s). \quad (15)$$

The leaking out of the spectral subspace is governed by the ‘off-diagonal’ part of Ω_{AD} . Namely, from estimate on $P\Omega_{AD}(s; \tau)P_\perp$ and $P_\perp\Omega_{AD}(s; \tau)P$ where $P = P(0)$ and $P_\perp = 1 - P$. This follows from

$$\begin{aligned} \|P_\tau(s) - P(s)\| &= \|P_\perp\Omega_{AD}(s; \tau)P - P\Omega_{AD}(s; \tau)P_\perp\| \\ &= \max\{\|P_\perp\Omega_{AD}(s; \tau)P\|, \|P\Omega_{AD}(s; \tau)P_\perp\|\} \end{aligned} \quad (16)$$

where

$$P_\tau(s) = U_\tau(s)PU_\tau^\dagger(s). \quad (17)$$

Ω_{AD} can be calculated via a Volterra-type equation:

$$\dot{\Omega}_{AD}(s; \tau) = K_{AD}(s)\Omega_{AD}(s; \tau) \tag{18}$$

where

$$K_{AD}(s) = -U_{AD}^\dagger(s; \tau)[\dot{P}, P](s)U_{AD}(s; \tau). \tag{19}$$

By standard arguments, the series

$$\Omega_{AD}(s; \tau) = \sum_0^\infty \Omega_i(s) \quad \Omega_0(s) = 1 \quad \Omega_{i+1}(s) = \int_0^s K_{AD}(t)\Omega_i(t) dt \tag{20}$$

is absolutely convergent[†].

It is not *a priori* clear that the series in equation (20) is an expansion in the small parameter of the adiabatic limit, $1/\tau$. To appreciate that it is, note the following lemma.

Lemma 1.

- (i) Ω_{2j} is diagonal in the sense that it maps Range P to Range P and Range P_\perp to Range P_\perp , at the same time Ω_{2j+1} is off-diagonal in the sense that it maps Range P to Range P_\perp and Range P_\perp to Range P . In particular, only the odd terms contribute to leaking.
- (ii) Let us denote $Q(s) := 1 - \Omega_{AD}(s)$, $f(\tau) := \sup_s \|Q(s)\|$. Then for all s ,

$$\|\Omega_{i+2}(s)\| < Cf(\tau) \sup_s \|\Omega_i(s)\|. \tag{21}$$

In particular

$$\begin{aligned} \|\Omega_j(s)\| &< Cf(\tau) & j = 1, 2 \\ \|\Omega_j(s)\| &< Cf^2(\tau) & j \geq 3. \end{aligned} \tag{22}$$

Remarks.

- (a) In the case where a gap condition holds, the uniform adiabatic theorem says that $f(\tau) = O(1/\tau)$. It follows that $\sup_s \|\Omega_{2i+1}(s)\| \leq \frac{C}{\tau^{i+1}}$. In the absence of a gap condition $f(\tau)$ may, in general, decay more slowly with τ .
- (b) For the lemma to be useful one needs a strong version of the uniform adiabatic theorem which guarantees that $Q(s)$ is small for large τ . This goes beyond the information given by the uniform adiabatic theorem quoted in the previous section. That is, the adiabatic evolution must approximate the true evolution both on Range P and its complement Range P_\perp .
- (c) Note that from the estimates above it follows, that when $\Omega_1(s) \gg f^2(\tau)$ for $s > 1$ that the leaky tails are determined by $\Omega_1(s)$.

Proof. The first part of the lemma is standard [2] and follows from the fact that $[\dot{P}, P]$ is off-diagonal. The second part follows from the identity

$$\Omega_{i+2}(s) = \int_0^s Q(t)K_{AD}(t)\Omega_i(t) dt - \int_0^s K_{AD}(t)\Omega_{AD}(t)Q^\dagger(t)\Omega_{i+1}(t) dt - Q(s)\Omega_{i+1}(s). \tag{23}$$

$K_{AD}(s)$ vanishes outside the interval $[0, 1]$, hence,

$$\begin{aligned} \|\Omega_{i+2}(s)\| &\leq f(\tau) \sup_s (\|K_{AD}(s)\| \cdot \|\Omega_i(s)\| + (\|K_{AD}(s)\| + 1)\|\Omega_{i+1}(s)\|) \\ &< 2f(\tau) (\sup_s \|K_{AD}(s)\| + 1) \sup_s \|\Omega_i(s)\|. \end{aligned} \tag{24}$$

□

[†] Recall that we assume $H(s)$ and $\dot{P}(s)$ are bounded.

4. Power tails in Friedrichs models

The Friedrichs model[†] is defined on the Hilbert space $\mathcal{H} = \mathbb{C} \oplus L^2(\mathbb{R}_+, d\mu(k))$ with $\inf(\text{support } \mu) \geq 0$ and $\mu(0) = 0$. A vector $\Psi \in \mathcal{H}$ is normalized by

$$\Psi = \begin{pmatrix} \omega \\ \psi(k) \end{pmatrix} \quad \|\Psi\|^2 = |\omega|^2 + \int_{\mathbb{R}_+} |\psi(k)|^2 d\mu(k) \quad \omega \in \mathbb{C}. \quad (25)$$

The Friedrichs Hamiltonian H acts on \mathcal{H} as

$$H\Psi = \begin{pmatrix} 0 & 0 \\ 0 & k \end{pmatrix} \begin{pmatrix} \omega \\ |\psi\rangle \end{pmatrix} = \begin{pmatrix} 0 \\ |k\psi\rangle \end{pmatrix}. \quad (26)$$

It has a ground state at threshold and the projection on the ground state P is

$$P = \begin{pmatrix} 1 & 0 \\ 0 & 0 \end{pmatrix}. \quad (27)$$

We consider the case when the time dependence is of the form $H(s) = V(s)HV^\dagger(s)$ with unitary $V(s)$ generated by

$$\dot{V}(s)V^\dagger(s) = [\dot{P}(s), P(s)] = i\dot{g}(s) \begin{pmatrix} 0 & \langle\varphi| \\ |\varphi\rangle & 0 \end{pmatrix} \quad (28)$$

where $0 \leq \dot{g}(s) \in C_0^\infty([0, 1])$, and

$$\int_0^x |\varphi|^2 d\mu(k) = O(x^{2\beta}) \quad \beta \geq 0 \quad (29)$$

for small x . Note that, due to the particular form of time dependence (equation (28)), one has that $H_{AD} = H_r$.

In the following proposition we borrow a result from [1].

Proposition 1. *For the Friedrichs model and the adiabatic evolution generated by H_{AD} , $f(\tau)$ of lemma 1 is such that*

$$f(\tau) \leq \begin{cases} O\left(\frac{1}{\tau}\right) & \beta > 1 \\ O\left(\frac{\log \tau}{\tau}\right) & \beta = 1 \\ O(\tau^{-\beta}) & \beta < 1 \end{cases} \quad (30)$$

for all s .

In the following proposition we present the main result of this section.

Proposition 2. *For the Friedrichs model the evolution of the state that starts as the bound state $\psi_\tau(0) \in \text{Range } P$, is such that for any $\epsilon > 0$*

$$C_1 \tau^{-\beta+\epsilon} > \text{dist}(\psi_\tau(s), \text{Range } P(s)) > C_2 \tau^{-\beta} \quad (31)$$

when $2 > \beta > 0$ and $s > 1$.

Proof. With the choice of H_{AD} all the even terms Ω_{2j} are diagonal while all the odd terms are off-diagonal. By lemma 1, $\Omega_j = O(f^2(\tau))$ for $j \geq 3$. We need to compute $\Omega_1(s)$ and,

[†] In some circles this is known as the Fanno model.

provided it dominates $f^2(\tau)$, we are done. Now, for $s > 1$,

$$\begin{aligned}
 P_{\perp}\Omega_1(s)P &= P_{\perp}\int_0^1 K(t) ds P \\
 &= i\begin{pmatrix} 0 & 0 \\ 0 & 1 \end{pmatrix}\int_0^1 \dot{g}(s)\exp\left(i\tau s\begin{pmatrix} 0 & 0 \\ 0 & k \end{pmatrix}\right)\begin{pmatrix} 0 & \langle\varphi| \\ |\varphi\rangle & 0 \end{pmatrix}\begin{pmatrix} 1 & 0 \\ 0 & 0 \end{pmatrix} ds \\
 &= i\int_0^1 \dot{g}(s)\begin{pmatrix} 0 & 0 \\ |\exp(i\tau sk)\varphi\rangle & 0 \end{pmatrix} ds \\
 &= i\begin{pmatrix} 0 & 0 \\ |\hat{g}(\tau k)\varphi\rangle & 0 \end{pmatrix} \tag{32}
 \end{aligned}$$

where \hat{g} stands for Fourier transform of g . Now, since g is positive, $|\hat{g}(k)|$ takes its maximal value at the origin. Since this function is continuous, for some a, b positive $|\hat{g}(k)| > b$ for $k \in [0, a]$. We can now estimate the norm of $P_{\perp}\Omega_1(s)P$ given below using equation (29):

$$\begin{aligned}
 \|P_{\perp}\Omega_1(s)P\|^2 &= \|\hat{g}(\tau k)\varphi(k)\|^2 = \int_0^{\infty} |\hat{g}(\tau k)|^2 |\varphi(k)|^2 d\mu(k) \\
 &> \int_0^a |\hat{g}(k)|^2 \left|\varphi\left(\frac{k}{\tau}\right)\right|^2 d\mu\left(\frac{k}{\tau}\right) \\
 &> b^2 \int_0^{\frac{a}{\tau}} |\varphi(k)|^2 d\mu(k) \\
 &= O\left(\frac{1}{\tau^{2\beta}}\right) \quad s > 1. \tag{33}
 \end{aligned}$$

Similarly,

$$\begin{aligned}
 \|P_{\perp}\Omega_1(s)P\|^2 &= \int_0^{\infty} |\hat{g}(\tau k)|^2 |\varphi(k)|^2 d\mu(k) \\
 &= \left\{ \int_0^{\frac{1}{\tau^{1-\epsilon}}} + \int_{\frac{1}{\tau^{1-\epsilon}}}^{\infty} \right\} |\hat{g}(\tau k)|^2 |\varphi(k)|^2 d\mu(k) \\
 &< C\left(\frac{1}{\tau^{2\beta(1-\epsilon)}} + |\hat{g}(\tau^{\epsilon})|^2\right) \quad s > 1. \tag{34}
 \end{aligned}$$

Since $|\hat{g}(k)|$ is decaying faster than any power at large k , the proof is complete. \square

Note that there is an improvement in the long-time limit over the uniform result for $2 > \beta \geq 1$.

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Appendix A. Slaved leaking

In this appendix we want to address an apparent puzzle associated with the (standard) long-time adiabatic theorem.

It is, in fact, surprising that a gap condition is all one needs for the fast decay in the long-time adiabatic theorem: suppose that $H(s)$ had one eigenvalue separated by a gap from the rest of the spectrum, which is purely absolutely continuous. For time $s \in [0, 1]$ the leaking to

the (instantaneous) absolutely continuous spectrum is $O(1/\tau)$. One would expect that once a piece leaks to the absolutely continuous part of the spectrum it would propagate to infinity, as states in the absolutely continuous spectrum invariably do. If this was the case, then it would be difficult to understand how the wave reconstructs itself so that, for times $s > 1$, the part in the absolutely continuous spectrum is smaller than any power in $1/\tau$. As we shall see from the proof of the adiabatic theorem to all orders below, the parts that leak to the absolutely continuous spectrum and is a finite power of τ is slaved to the instantaneous eigenvalue, and disappears for time $s > 1$. It does not propagate. Only the terms smaller than any power in $1/\tau$ are free to propagate to infinity.

The following lemma is taken, verbatim, from [18].

Lemma 2. Let $R(s, z)$ be the resolvent of $H(s)$ and define the tilde operation by

$$\tilde{X}(s) = -\frac{1}{2\pi i} \int_{\Gamma} R(s, z) X(s) R(s, z) dz \quad (\text{A.1})$$

where Γ is a contour enclosing the part of the spectrum associated with $P(s)$. Let, $X(s)$ and $Y(s)$ be bounded, then

$$P_{\perp} \int_0^s U_{AD}^{\dagger}(t) X(t) U_{AD} P Y(t) dt = \frac{i}{\tau} P_{\perp} \left(U_{AD}^{\dagger}(t) \tilde{X}(t) U_{AD} P Y(t) \Big|_0^s - \int_0^s U_{AD}^{\dagger}(t) \dot{\tilde{X}}(t) U_{AD} P Y(t) dt - \int_0^s U_{AD}^{\dagger}(t) \tilde{X}(t) U_{AD} P \dot{Y}(t) dt \right). \quad (\text{A.2})$$

If we now apply the lemma with $X(s) = [\dot{P}, P](s)$ to Ω_{j+1} we get

$$P_{\perp} \Omega_{j+1}(s) = P_{\perp} \int_0^s U_{AD}^{\dagger}(t) X(t) U_{AD} P \Omega_j(t) dt = -\frac{i}{\tau} P_{\perp} \left(U_{AD}^{\dagger}(s) \tilde{X}(s) U_{AD}(s) P \Omega_j(s) - \int_0^s U_{AD}^{\dagger}(t) \dot{\tilde{X}}(t) U_{AD} P \Omega_j(t) dt - \int_0^s U_{AD}^{\dagger}(t) \tilde{X}(t) U_{AD} P K(t) \Omega_{j-1}(t) dt \right). \quad (\text{A.3})$$

Each of the *integrals* in the above expression are once again of the form such that the lemma can be applied again and again. We see that the power terms that occur for $0 < s < 1$ are boundary terms, proportional to $\dot{P}(s)$, and instantaneously reflect what happens to the bound state. That is, the power tails in $\text{Range } P_{\perp}(s)$ are slaved to the state in $\text{Range } P(s)$.

Appendix B. Asymptotics of Fourier transforms in C_0^{∞}

When $g \in C_0^{\infty}$ its Fourier transform decays faster than any power, but not quite exponentially. The canonical example of such a function is

$$g(s) = \exp\left(\frac{-1}{(1-s)(1+s)}\right) \quad s \in [-1, 1] \quad (\text{B.1})$$

and zero otherwise. It is of some interest to have an explicit asymptotic expansion of its Fourier transform.

Lemma 3. $\hat{g}(p)$ is an entire even function of p , and on the real p -axis it has the asymptotic behaviour

$$\hat{g}(p) = 2 \frac{\sqrt{\pi}}{(2e)^{1/4}} \cdot \frac{\exp(-\sqrt{p})}{p^{3/4}} \cdot \left(\cos\left(p - \sqrt{p} - \frac{3}{8}\pi\right) + O(1/\sqrt{p}) \right). \quad (\text{B.2})$$

Proof. The fact that the function is even and entire follows directly from the definition

$$\hat{g}(p) = \int_{-1}^1 \exp(ips) \cdot \exp\left(\frac{-1}{1-s^2}\right) ds = \int_{-1}^1 \cos(ps) \cdot \exp\left(\frac{-1}{1-s^2}\right) ds. \quad (\text{B.3})$$

The asymptotics follow from the Laplace saddle point method. Let

$$h(s, p) = -\frac{1}{1-s^2} + isp \quad (\text{B.4})$$

denote the logarithm of the integrand. The integral is estimated by Gaussian integrals near the appropriate extrema of h . The extremum at s_0 contributes

$$\sqrt{\frac{2\pi}{-h''(s_0, p)}} \cdot \exp h(s_0, p) \quad (\text{B.5})$$

where the prime denotes a partial derivative with respect to s .

The extrema are the solutions of the quartic equation

$$h'(s, p) = -\frac{2s}{(1-s^2)^2} + ip = 0. \quad (\text{B.6})$$

When p is large, two of the solutions are close to 1 and the other two are close to -1 . To leading order in p the two extrema near 1 are

$$s_0(p) = (1 \pm \delta(p)) \quad \delta(p) = \frac{1}{\sqrt{2ip}} + O\left(\frac{1}{p^{3/2}}\right) \quad (\text{B.7})$$

where

$$h(1 \pm \delta, p) = ip \pm (1+i)\sqrt{p} - \frac{1}{4} + O(\delta). \quad (\text{B.8})$$

It is now clear that the right way to deform the contour is so that it goes through $1 - \delta$, and the saddle point at $1 + \delta$ should be avoided. Now, to leading order in p , the second derivative there is

$$\begin{aligned} h''(s_0, p) &= -\frac{8s^2}{(1-s^2)^3} + O(p) \\ &= -(2ip)^{3/2}(1 + O(\delta)) \\ &= -(2p)^{3/2} \exp\left(\frac{3\pi i}{4}\right) \cdot (1 + O(\delta)). \end{aligned} \quad (\text{B.9})$$

Substituting in equation (B.5) one finds the contribution from the saddle points near 1 to be

$$\frac{\sqrt{\pi}}{(2e)^{1/4}} \cdot \frac{\exp(-\sqrt{p})}{p^{3/4}} \cdot \exp(i(p - \sqrt{p} - 3\pi/8))(1 + O(1/\sqrt{p})). \quad (\text{B.10})$$

Using the fact that $\hat{g}(p)$ is real valued for real p , the saddle near -1 must give the complex conjugate of this, and the result follows. \square

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