

A stable, polynomial-time algorithm for the eigenpair problem

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Abstract. We describe algorithms for computing eigenpairs (eigenvalue-eigenvector pairs) of a complex $n \times n$ matrix A . These algorithms are numerically stable, strongly accurate, and theoretically efficient (i.e., polynomial-time). We do not believe they outperform in practice the algorithms currently used for this computational problem. The merit of our paper is to give a positive answer to a long-standing open problem in numerical linear algebra.

So the problem of devising an algorithm [for the eigenvalue problem] that is numerically stable and globally (and quickly!) convergent remains open.

J. Demmel [25, page 139]

*Partially supported by Agencia Nacional de Investigación e Innovación (ANII), Uruguay, and by CSIC group 618

†partially supported by the research projects MTM2010-16051 and MTM2014-57590 from Spanish Ministry of Science MICINN

‡Partially funded by DFG research grant BU 1371/2-2

§Partially funded by a GRF grant from the Research Grants Council of the Hong Kong SAR (project number CityU 100813).

1 Introduction

1.1 The problem

The quotation from Demmel opening this article, though possibly puzzling for those who day-to-day satisfactorily solve eigenvalue problems, summarizes a long-standing open problem in numerical linear algebra. The first algorithm that comes to mind for computing eigenvalues —to compute the characteristic polynomial χ_A of A and then compute (i.e., approximate) its zeros— has proved to be numerically unstable. The so called Wilkinson's polynomial,

$$w(x) := \prod_{i=1}^{20} (x - i) = x^{20} + w_{19}x^{19} + \cdots + w_1x + w_0$$

is often used to illustrate this fact. For a diagonal matrix D with diagonal entries $1, 2, \dots, 20$ (and therefore with $\chi_D(x) = w(x)$) an error of 2^{-23} in the computation of $w_{19} = -210$ produces, even if the rest of the computation is done error-free, catastrophic variations in the zeros of χ_D . For instance, the zeros at 18 and 19 collide into a double zero close to 18.62, which will unfold into two complex conjugate zeros if the error in w_{19} is further increased. And yet, there is nothing wrong in the nature of D (in numerical analysis terms, as we will see below, D is a well-conditioned matrix for the eigenvalue problem). The trouble appears to lie in the method.

Barred from using this immediate algorithm due to its numerical instability, researchers devoted efforts to come up with alternate methods which would appear to be stable. Among those proposed, the one that is today's algorithm of choice is the iterated QR with shifts. This procedure behaves quite efficiently in general and yet, as Demmel pointed out in 1997 [25, p. 139],

after more than 30 years of dependable service, convergence failures of this algorithm have quite recently been observed, analyzed, and patched [...]. But there is still no global convergence proof, even though the current algorithm is considered quite reliable.

Our initial quotation followed these words in Demmel's text. It asked for an algorithm which will be numerically stable and for which, convergence, and if possible small complexity bounds, can be established. Today, 17 years after Demmel's text, this demand retains all of its urgency: it is not known if any of the standard numerical linear algebra algorithms satisfies the properties above. For example:

- The unshifted QR algorithm terminates with probability 1 but is probably infinite average cost if approximations to the eigenvectors are to be output (see [32]).
- The QR algorithm with Rayleigh Quotient shift fails for open sets of real input matrices (see [8, 9]).

- We do not know whether the Francis (double) shift algorithm converges generally on real or complex matrices, nor an estimate of its average cost.
- Other algorithms in modern texts are analyzed but don't estimate the (necessarily infinite in the worst case) number of iterations, which usually relies on experimental results, see for example [40] which uses a divide and conquer algorithm or [26] which is on turn based on the algorithm in [6].

Algorithms which output approximate eigenvalues without accompanying approximate eigenvectors might be easier to analyze. The experimental evidence of [41] for symmetric matrices suggests that many of the algorithms in use are of average finite cost and even that there is some universality. An informal explanation of this fact is that the eigenvalues of symmetric matrices are very well conditioned, see for example [56, eq. (1.5)]. But eigenvectors are another matter. When the matrices are close to having multiple eigenvalues the condition of the eigenvector tends to infinity. For example, even for 2×2 symmetric matrices, any pair of orthogonal vectors (a, b) and $(-b, a)$ are the eigenvectors of a matrix

$$\begin{pmatrix} 1 + \varepsilon_1 & \varepsilon_3 \\ \varepsilon_3 & 1 + \varepsilon_2 \end{pmatrix}$$

for $|\varepsilon_i|, i = 1, 2, 3$, arbitrarily small.

The only goal of this paper is to give a positive answer to Demmel's question. The set of our main results can be informally stated as follows.

Main results We exhibit algorithms which on input a complex matrix A with complex Gaussian entries generate (with probability 1) an “excellent” approximation to one of (or all) the (eigenvalue, eigenvector) pairs of A . Some of these algorithms are deterministic while some are randomized. Their running time (expected running time for the randomized case) is polynomial in n on average (w.r.t. A) as well as under a standard smoothed analysis.

More precisely, the average complexity bounds we prove, for $n \times n$ matrices, are $\mathcal{O}(n^7)$ —for the computation of a single eigenpair with either a deterministic or a randomized algorithm— and $\mathcal{O}(n^9)$ —for the computation of all eigenpairs with a deterministic algorithm. Note that these are just upper bounds: the practical performance of the algorithms might be better (see §2.10). The precise statements of the main results are in Theorems 2.25, 2.28, and 2.35.

1.2 A few words on approximations

It must be said upfront that we do not think the algorithm we propose will outperform, in general, iterated QR with shifts. It nonetheless possesses some worthy features which we want to describe in this introduction. The key one, we already mentioned, is that both convergence and complexity bounds can be established for it. It is also numerically stable. But in addition, it is *strongly accurate*.

A starting point to understand the meaning of this last claim, is the observation that there are two different obstructions to the exact computation of an eigenpair. Firstly, the use of finite precision, and the ensuing errors accumulating during the computational process. The expression *numerically stable* is usually vested on algorithms for which this accumulated error on the computed quantities is not much larger than that resulting from an error-free computation on an input datum which has been read (and approximated) with machine precision. Secondly, the nonlinear character of the equations defining eigenvalues and eigenvectors in terms of the given matrix. For $n \geq 5$, we learned from Abel and Galois, we cannot write down these eigenvalues in terms of the matrix' entries, not even using radicals, and the same remains true for eigenvectors. Hence, we can only compute approximations of them and this is so *even assuming infinite precision in the computation*.

The expression *strongly accurate* refers to the quality of these approximations. It is common to find in the literature (at least) three notions of approximation which we next briefly describe. To simplify, we illustrate with the computation of a value $\zeta \in \mathbb{C}$ from a datum $A \in \mathbb{C}^N$ (and the reader may well suppose that this computation is done with infinite precision). We let $\tilde{\zeta}$ be the quantity actually computed and we consider the following three requirements on it:

Backward approximation. The element $\tilde{\zeta}$ is the solution of a datum \tilde{A} close to A .

Given $\varepsilon > 0$, we say that $\tilde{\zeta}$ is an ε -backward approximation when $\|A - \tilde{A}\| \leq \varepsilon$ (resp. $\|A - \tilde{A}\| \leq \varepsilon \|A\|$ if we are interested in relative errors).

Forward approximation. The quantity $\tilde{\zeta}$ is close to ζ . Given $\varepsilon > 0$, we say that $\tilde{\zeta}$ is an ε -forward approximation when $|\zeta - \tilde{\zeta}| \leq \varepsilon$ (resp. $|\zeta - \tilde{\zeta}| \leq \varepsilon |\zeta|$).

Approximation à la Smale. An appropriate version of Newton's iteration, starting at $\tilde{\zeta}$, converges immediately, quadratically fast, to ζ , either in absolute or relative error.

These requirements are often, but not always, increasingly demanding. For instance, if ζ is an ε -backward approximation then the forward error $|\zeta - \tilde{\zeta}|$ is bounded, roughly, by $\varepsilon \text{cond}(A)$. Here $\text{cond}(A)$ is the condition number of A , a quantity usually greater than 1. So, in general, ε -backward approximations are not ε -forward approximations, and if A is poorly conditioned $\tilde{\zeta}$ may be a poor forward approximation of ζ (we note, however, that backward approximations do not necessarily exist, and that condition numbers smaller than 1 do occur). We also observe that if $\tilde{\zeta}$ is an approximation à la Smale we can obtain an ε -forward approximation from $\tilde{\zeta}$ by performing $\mathcal{O}(\log |\log \varepsilon|)$ Newton's steps (see for example Theorem 1.1 below). To obtain an approximation à la Smale from an ε -forward approximation we need, in contrast, that ε will be of the order of the inverse of the condition of the zero (see Theorem 2.12 below).

When we say that our algorithm is strongly accurate, we refer to the fact that the returned eigenpairs are approximations à la Smale of true eigenpairs.

In our case, we can not only efficiently compute ε -forward approximations as above but also with respect to relative error. These are pairs (ζ, w) satisfying

$$d_{\mathbb{P}}(w, v) \leq \varepsilon \quad \text{and} \quad |\zeta - \lambda| \leq \varepsilon |\lambda|$$

for some true eigenpair (λ, v) of A . Here $d_{\mathbb{P}}(w, v)$ denotes the angle between w and v (note that here the scaling of eigenvectors renders the relativization of the error moot).

Theorem 1.1 *We exhibit an algorithm that, given a matrix $A \in \mathbb{C}^{n \times n}$, an approximate eigenpair (ζ, w) returned by any of the algorithms in the main results, and an accuracy $\varepsilon \in (0, 1/2)$, produces an ε -forward approximation (in relative error) of the approximated true eigenpair (λ, v) . The algorithm terminates provided $\lambda \neq 0$. Its average cost over Gaussian matrices A (independently of the chosen approximate eigenpair) is $\mathcal{O}(n^3 \log_2 \log_2(n/\varepsilon))$.*

Combining our main results with Theorem 1.1 we can thus compute ε -forward approximations (in relative error) of all eigenpairs of random Gaussian matrices with average running time $\mathcal{O}(n^9 + n^3 \log_2 \log_2(n/\varepsilon))$. See §11 for a proof of Theorem 1.1.

1.3 A few words on complexity

The cost, understood as the number of arithmetic operations performed, of computing an approximation of an eigenpair for a matrix $A \in \mathbb{C}^{n \times n}$, depends on the matrix A itself. Actually, and this is a common feature in numerical analysis, it depends on the condition $\text{cond}(A)$ of the matrix A . But this condition is not known a priori. It was therefore advocated by Smale [51] to eliminate this dependency in complexity bounds by endowing data space with a probability distribution and estimating expected costs. This idea has its roots in early work of Goldstine and von Neumann [57].

In our case, data space is $\mathbb{C}^{n \times n}$, and a common probability measure to endow it with is the standard Gaussian. Expectations of cost w.r.t. this measure yield expressions in n usually referred to as *average cost*. A number of considerations, including the suspicion that the use of the standard Gaussian could result in complexity bounds which are too optimistic compared with “real life”, prompted Spielman and Teng to introduce a different form of probabilistic analysis, called *smoothed analysis*. In this, one replaces the average analysis’ goal of showing that

for a random A it is unlikely that the cost for A will be large
by the following

for all \hat{A} , it is unlikely that a slight random perturbation $A = \hat{A} + \Delta A$
will require a large cost.

The expectations obtained for a smoothed analysis will now be functions of both the dimension n and some measure of dispersion for the random perturbations (e.g., a variance).

Smoothed analysis was first used for the simplex method of linear programming [54]. Some survey expositions of its rationale are in [53, 55, 17]. One may argue that it has been well accepted by the scientific community from the fact that Spielman and Teng were awarded the Gödel 2008 and Fulkerson 2009 prizes for it (the former by the theoretical computer science community and the latter by the optimization community). Also, in 2010, Spielman was awarded the Nevanlinna prize, and smoothed analysis appears in the laudatio of his work.

In this paper we will exhibit bounds for the cost of our algorithm both for average and smoothed analyses.

1.4 A few words on numerical stability

The algorithm we deal with in this paper belongs to the class of homotopy continuation methods. Experience has shown that algorithms in this class are very stable and stability analyses have been done for some of them, e.g. [16, 11, 23]. Because of this, we will assume infinite precision all along this paper and steer clear of any form of stability analysis. We nonetheless observe that such an analysis can be easily carried out following the steps in the papers mentioned above.

1.5 Previous and related work

Homotopy continuation methods go back, at least, to the work of Lahaye [33]. A detailed survey of their use to solve polynomial equations is in [35]. More explicit focus in eigenvalue computations is considered in [21, 36, 37, 38] but we do not know of any serious attempt to implement them.

In the early 1990s Shub and Smale set up a program to understand the cost of solving square systems of complex polynomial equations using homotopy methods. In a collection of articles [45, 46, 47, 48, 49], known as the *Bézout series*, they put in place many of the notions and techniques that occur in this article. The Bézout series did not, however, conclusively settle the understanding of the cost above, and in 1998 Smale proposed it as the 17th in his list of problems for the mathematicians of the 21st century [52]. A probabilistic solution to this problem was found in [12, 13], then a deterministic quasipolynomial solution was described in [18]. Finally, a deterministic polynomial solution was recently found by Lairez in [34] and is now considered fully answered by the community.

The results in these papers cannot be directly used for the eigenpair problem since instances of the latter are ill-posed as polynomial systems. But the intervening ideas can be reshaped to attempt a tailor-made analysis for the eigenpair problem. A major step in this direction was done by Armentano in his PhD thesis (see [3] and its precedent [24]), where the condition number μ for the eigenpair problem was

exhaustively studied. A further step was taken in [5] where μ was used to analyze a randomized algorithm for the Hermitian eigenpair problem.

Our paper follows this stream of research.

Since the appearance of our work, Paul Breiding has successfully generalized the ideas of our paper to the computation of eigenpairs of homogeneous polynomial systems, see [15].

1.6 Structure of the exposition

The remainder of this paper is divided into two parts. In the first one, §2 below, we introduce all the technical preliminaries, we describe with details the algorithms, and we state our main results (Theorems 2.25, 2.28, and 2.35). The condition number μ , Newton's method, the notion of approximate eigenpair, and Gaussian distributions are among these technical preliminaries. The second part, which occupies us in the subsequent sections, is devoted to proofs. Some short proofs are included in §2 as well.

2 Preliminaries, Basic Ideas, and Main Result

2.1 Spaces and Metric Structures

Let $\mathbb{C}^{n \times n}$ be the set of $n \times n$ complex matrices. We endow this complex linear space with the *Frobenius Hermitian product* $\langle \cdot, \cdot \rangle_F$ and the associated *Frobenius norm* $\|\cdot\|_F$ given by

$$\langle A, B \rangle_F := \text{trace} (B^* A) = \sum_{i,j=1}^n a_{ij} \overline{b_{ij}}, \quad \|A\|_F = \langle A, A \rangle_F^{1/2},$$

where $A = (a_{ij})$ and $B = (b_{ij})$. The unit sphere will be denoted by $\mathbb{S}(\mathbb{C}^{n \times n})$ or simply by \mathbb{S} . We endow the product vector space $\mathbb{C}^{n \times n} \times \mathbb{C}$ with the canonical Hermitian inner product structure and its associated norm structure and (Euclidean) distance.

The space \mathbb{C}^n is equipped with the canonical Hermitian inner product $\langle \cdot, \cdot \rangle$. We denote by $\mathbb{P}(\mathbb{C}^n)$ its associated projective space. This is a smooth manifold which carries a natural Riemannian metric, namely, the real part of the *Fubini-Study metric* on $\mathbb{P}(\mathbb{C}^n)$. The Fubini-Study metric is the Hermitian structure on $\mathbb{P}(\mathbb{C}^n)$ given in the following way: for $x \in \mathbb{C}^n$,

$$\langle w, w' \rangle_x := \frac{\langle w, w' \rangle}{\|x\|^2}, \quad (1)$$

for all w, w' in the Hermitian complement $x^\perp = \{v \in \mathbb{C}^n \mid \langle x, v \rangle = 0\}$ of x in \mathbb{C}^n . We denote by $d_{\mathbb{P}}$ the associated Riemannian distance. An explicit formula for that

distance (see for example [14, p. 226]) is

$$d_{\mathbb{P}}(v, w) = \arccos \frac{|\langle v, w \rangle|}{\|v\| \cdot \|w\|}. \quad (2)$$

Note that this formula makes sense for $v, w \in \mathbb{C}^n$ (as the distance between the respective projective classes).

The space $\mathbb{C}^{n \times n} \times \mathbb{C} \times \mathbb{P}(\mathbb{C}^n)$ is endowed with the Riemannian product structure. The resulting distance equals

$$\text{dist}((A, \lambda, v), (A', \lambda', v'))^2 := \|A - A'\|_F^2 + |\lambda - \lambda'|^2 + d_{\mathbb{P}}(v, v')^2.$$

We will only use this distance on $\mathbb{S} \times \mathbb{C} \times \mathbb{P}(\mathbb{C}^n)$. Note that for $A, A' \in \mathbb{S}$, the distance $\text{dist}((A, \lambda, v), (A', \lambda', v'))$ is smaller than or equal to the natural geodesic (product) distance in $\mathbb{S} \times \mathbb{C} \times \mathbb{P}(\mathbb{C}^n)$. For any nonzero matrix $A \in \mathbb{C}^{n \times n}$ (not necessarily of unit norm) we will write

$$\text{dist}_A((\lambda, v), (\lambda', v'))^2 := \frac{|\lambda - \lambda'|^2}{\|A\|_F^2} + d_{\mathbb{P}}(v, v')^2.$$

Note that for any nonzero $A \in \mathbb{C}^{n \times n}$, dist_A is a distance function in $\mathbb{C} \times \mathbb{P}(\mathbb{C}^n)$ and if $A \in \mathbb{S}$, then $\text{dist}_A((\lambda, v), (\lambda', v')) = \text{dist}((A, \lambda, v), (A, \lambda', v'))$.

2.2 The Varieties \mathcal{V} , \mathcal{W} , Σ' and Σ

We define the *solution variety* for the eigenpair problem as

$$\mathcal{V} = \mathcal{V}_n := \{(A, \lambda, v) \in \mathbb{C}^{n \times n} \times \mathbb{C} \times \mathbb{P}(\mathbb{C}^n) : (A - \lambda \text{Id})v = 0\}.$$

Proposition 2.1 *The solution variety \mathcal{V} is a smooth submanifold of $\mathbb{C}^{n \times n} \times \mathbb{C} \times \mathbb{P}(\mathbb{C}^n)$, of the same dimension as $\mathbb{C}^{n \times n}$.*

PROOF. See [3, Proposition 2.2]. □

The set \mathcal{V} inherits the Riemannian structure of the ambient space. Associated to it there are natural projections:

$$\begin{array}{ccc} & \mathcal{V} & \\ \pi_1 \swarrow & & \searrow \pi_2 \\ \mathbb{C}^{n \times n} & & \mathbb{C} \times \mathbb{P}(\mathbb{C}^n). \end{array} \quad (3)$$

Because of Proposition 2.1, the derivative $D\pi_1$ at (A, λ, v) is a linear operator between spaces of equal dimension. The *subvariety \mathcal{W} of well-posed triples* is the

subset of triples $(A, \lambda, v) \in \mathcal{V}$ for which $D\pi_1(A, \lambda, v)$ is an isomorphism. In particular, when $(A, \lambda, v) \in \mathcal{W}$, the projection π_1 has a branch of its inverse (locally defined) taking $A \in \mathbb{C}^{n \times n}$ to $(A, \lambda, v) \in \mathcal{V}$.

Let $P_{v^\perp} : \mathbb{C}^n \rightarrow v^\perp$ be the orthogonal projection. Given $(A, \lambda, v) \in \mathbb{C}^{n \times n} \times \mathbb{C} \times \mathbb{P}(\mathbb{C}^n)$, we let $A_{\lambda, v} : v^\perp \rightarrow v^\perp$ be the linear operator given by

$$A_{\lambda, v} := P_{v^\perp} \circ (A - \lambda \text{Id})|_{v^\perp}. \quad (4)$$

If we choose a representative such that $\|v\| = 1$ and we assume that $A_{\lambda, v}$ is invertible, then we have

$$i_{\mathbb{C}^n} A_{\lambda, v} P_{v^\perp} = (\text{Id} - vv^*)(A - \lambda \text{Id})(\text{Id} - vv^*), \quad (5)$$

and

$$i_{\mathbb{C}^n} A_{\lambda, v}^{-1} P_{v^\perp} = ((\text{Id} - vv^*)(A - \lambda \text{Id})(\text{Id} - vv^*))^\dagger, \quad (6)$$

where $i_{\mathbb{C}^n} : v^\perp \rightarrow \mathbb{C}^n$ is the inclusion and † denotes Moore–Penrose pseudoinverse.

The set of well-posed triples is exactly

$$\mathcal{W} = \{(A, \lambda, v) \in \mathcal{V} : A_{\lambda, v} \text{ is invertible}\}, \quad (7)$$

see [3, Lemma 2.7]). We define $\Sigma' := \mathcal{V} \setminus \mathcal{W}$ to be the variety of *ill-posed triples*, and $\Sigma = \pi_1(\Sigma') \subset \mathbb{C}^{n \times n}$ the *discriminant variety*, i.e., the subset of *ill-posed inputs*.

Remark 2.2 From (7) it is clear that the subset Σ' is the set of triples $(A, \lambda, v) \in \mathcal{V}$ such that λ is an eigenvalue of A of algebraic multiplicity at least 2. Hence Σ is the set of matrices $A \in \mathbb{C}^{n \times n}$ with multiple eigenvalues, and for $A \in \mathbb{C}^{n \times n} \setminus \Sigma$, the eigenvalues of A are pairwise different and $\pi_1^{-1}(A)$ is the set of triples $(A, \lambda_1, v_1), \dots, (A, \lambda_n, v_n)$, where (λ_i, v_i) , $i = 1, \dots, n$, are the eigenpairs of A .

Proposition 2.3 *The discriminant variety $\Sigma \subset \mathbb{C}^{n \times n}$ is a complex algebraic hypersurface. Consequently, for all $n \geq 2$, we have $\dim_{\mathbb{R}} \Sigma = 2n^2 - 2$.*

PROOF. See [19, Proposition 20.18]. \square

2.3 Unitary invariance

Let \mathcal{U}_n be the group of $n \times n$ unitary matrices. The group \mathcal{U}_n naturally acts on $\mathbb{P}(\mathbb{C}^n)$ by $U \cdot w := Uw$. In addition, \mathcal{U}_n acts on $\mathbb{C}^{n \times n}$ by conjugation (i.e., $U \cdot A := UAU^{-1}$), and on $\mathbb{C}^{n \times n} \times \mathbb{C}$ by $U \cdot (A, \lambda) := (UAU^{-1}, \lambda)$. These actions define an action on the product space $\mathbb{C}^{n \times n} \times \mathbb{C} \times \mathbb{P}(\mathbb{C}^n)$, namely,

$$U \cdot (A, \lambda, v) := (UAU^{-1}, \lambda, Uv). \quad (8)$$

Remark 2.4 The varieties \mathcal{V} , \mathcal{W} , Σ' , and Σ , are invariant under the action of \mathcal{U}_n (see [3] for details).

2.4 Condition of a triple

In a nutshell, condition numbers measure the worst possible output error resulting from a small perturbation on the input data. More formally, a condition number is the operator norm of the derivative of a solution map such as the branches of π_1^{-1} mentioned in §2.1 above, (see [19, §14.1.2] for a general exposition).

In the case of the eigenpair problem, one can define two condition numbers for the eigenvalue and the eigenvector, respectively, and formulas for both of them have been known at least since [56]. Armentano has shown that one can merge the two in a single one (see §3 in [3] for details). Deviating slightly from [3], we define the *condition number* of $(A, \lambda, v) \in \mathbb{C}^{n \times n} \times \mathbb{C} \times \mathbb{C}^n$ as

$$\mu(A, \lambda, v) := \|A\|_F \|A_{\lambda, v}^{-1}\|, \quad (9)$$

(or ∞ if $A_{\lambda, v}$ is not invertible) where $\|\cdot\|$ is the operator norm. This coincides with $\mu_v(A, \lambda, v)$ in [3]. Note that from (6), if (A, λ, v) is such that $\mu(A, \lambda, v) < \infty$, and $\|v\| = 1$ then:

$$\mu(A, \lambda, v) = \|A\|_F \left\| ((\mathbf{Id} - vv^*)(A - \lambda \mathbf{Id})(\mathbf{Id} - vv^*))^\dagger \right\|. \quad (10)$$

Remark 2.5 The condition number μ is invariant under the action of the unitary group \mathcal{U}_n , i.e., $\mu(UAU^{-1}, \lambda, UV) = \mu(A, \lambda, v)$ for all $U \in \mathcal{U}_n$, and scale invariant on the first two components, i.e., $\mu(sA, s\lambda, v) = \mu(A, \lambda, v)$ for all $s \in \mathbb{C} \setminus \{0\}$.

Lemma 2.6 (Lemma 3.8 in [3]) *For $(A, \lambda, v) \in \mathcal{V}$ we have $\mu(A, \lambda, v) \geq \frac{1}{\sqrt{2}}$. \square*

The essence of condition numbers is that they measure how much may outputs vary when inputs are slightly perturbed. The following result, which we will prove in §3, quantifies this property for μ .

Proposition 2.7 *Let $\Gamma : [0, 1] \rightarrow \mathcal{V}$, $\Gamma(t) = (A_t, \lambda_t, v_t)$ be a smooth curve such that A_t lies in the unit sphere of $\mathbb{C}^{n \times n}$, for all t . Write $\mu_t := \mu(\Gamma(t))$. Then we have, for all $t \in [0, 1]$,*

$$|\dot{\lambda}_t| \leq \sqrt{1 + \mu_t^2} \|\dot{A}_t\|, \quad \|\dot{v}_t\| \leq \mu_t \|\dot{A}_t\|.$$

In particular,

$$\|\dot{\Gamma}(t)\| \leq \sqrt{6} \mu_t \|\dot{A}_t\|.$$

Condition numbers are generally associated to input data. In the case of a problem with many possible solutions (of which returning an eigenpair of a given matrix is a clear case) one can derive the condition of a data from a notion of condition for each of these solutions. A discussion of this issue is given in [19, §6.8]. For the purposes of this paper, we will be interested in two such derivations: the *maximum condition number* of A ,

$$\mu_{\max}(A) := \max_{1 \leq j \leq n} \mu(A, \lambda_j, v_j),$$

and the *mean square condition number* of A ,

$$\mu_{\mathbf{av}}(A) := \left(\frac{1}{n} \sum_{j=1}^n \mu^2(A, \lambda_j, v_j) \right)^{\frac{1}{2}} = \left(\frac{1}{n} \sum_{j=1}^n \|A\|_F^2 \|A_{\lambda_j, v_j}^{-1}\|^2 \right)^{\frac{1}{2}}.$$

Condition numbers themselves vary in a controlled manner. The following Lipschitz property and its corollary make this statement precise.

Theorem 2.8 *Let $A, A' \in \mathbb{S}$, let $v, v' \in \mathbb{C}^n$ be nonzero, and let $\lambda, \lambda' \in \mathbb{C}$. Suppose that*

$$\mu(A, \lambda, v) \text{dist}((A, \lambda, v), (A', \lambda', v')) \leq \frac{\varepsilon}{4\sqrt{3}} \quad \text{for some } \varepsilon \in (0, 1).$$

Then we have

$$\frac{1}{1+\varepsilon} \mu(A, \lambda, v) \leq \mu(A', \lambda', v') \leq \frac{1}{1-\varepsilon} \mu(A, \lambda, v).$$

Corollary 2.9 *Let $A \in \mathbb{S}$, $A \notin \Sigma$, and let $A' \in \mathbb{S}$ be such that*

$$\|A - A'\|_F \leq \frac{\varepsilon}{50\mu_{\max}^2(A)}, \quad \text{for some } \varepsilon \in (0, 1/2).$$

Then, $A' \notin \Sigma$ and $(1+\varepsilon)^{-1}\mu_{\max}(A) \leq \mu_{\max}(A') \leq (1-\varepsilon)^{-1}\mu_{\max}(A)$.

We give the proofs of Theorem 2.8 and Corollary 2.9 in §4 below.

We close this paragraph observing that restricted to the class of normal matrices, the condition number μ admits the following elegant characterization.

Lemma 2.10 (Lemma 3.12 in [3]) *Let $A \in \mathbb{C}^{n \times n} \setminus \Sigma$ be normal, and let $(\lambda_1, v_1), \dots, (\lambda_n, v_n)$ be its eigenpairs. Then*

$$\mu(A, \lambda_1, v_1) = \frac{\|A\|_F}{\min_{2 \leq j \leq n} |\lambda_j - \lambda_1|}.$$

□

2.5 Newton's method and approximate eigenpairs

For a nonzero matrix $A \in \mathbb{C}^{n \times n}$, we define the *Newton map* associated to A ,

$$N_A : \mathbb{C} \times (\mathbb{C}^n \setminus \{0\}) \rightarrow \mathbb{C} \times (\mathbb{C}^n \setminus \{0\}),$$

by

$$N_A \begin{pmatrix} \lambda \\ v \end{pmatrix} = \begin{pmatrix} \lambda \\ v \end{pmatrix} - \begin{pmatrix} \dot{\lambda} \\ \dot{v} \end{pmatrix}, \quad \text{where} \quad \begin{pmatrix} \dot{\lambda} \\ \dot{v} \end{pmatrix} = (DF_A(\lambda, v) |_{\mathbb{C} \times v^\perp})^{-1} F_A \begin{pmatrix} \lambda \\ v \end{pmatrix}$$

and $F_A(\lambda, v) = (A - \lambda \mathbf{Id})v$ is the evaluation map. This is a rational map (it is only defined on an open subset of $\mathbb{C} \times (\mathbb{C}^n \setminus \{0\})$). It was introduced in [3] as a Newton-like operator associated to the evaluation map F_A , and the following formulas were obtained for \dot{v} and $\dot{\lambda}$ (recall the definition of $A_{\lambda, v}$ from (4)):

$$\dot{v} = i_{\mathbb{C}^n} A_{\lambda, v}^{-1} P_{v^\perp} A v, \quad \dot{\lambda} = \frac{\langle (\lambda \mathbf{Id} - A)(v - \dot{v}), v \rangle}{\langle v, v \rangle}. \quad (11)$$

The map N_A is defined for every $(\lambda, v) \in \mathbb{C} \times (\mathbb{C}^n \setminus \{0\})$ such that $A_{\lambda, v}$ is invertible. It is immediate to check that for $k \geq 0$ and $z \in \mathbb{C}$ we have

$$N_{zA}^k \begin{pmatrix} z\lambda \\ v \end{pmatrix} = \begin{pmatrix} z & 0 \\ 0 & 1 \end{pmatrix} N_A^k \begin{pmatrix} \lambda \\ v \end{pmatrix}, \quad (12)$$

where the superindex k means k iterations. See [3, Sec. 4] for more details.

The notion of approximate solution as a point where Newton's method converges to a true solution immediately and quadratically fast was introduced by Steve Smale [50]. It allows to elegantly talk about approximations without dependencies on pre-established accuracies. In addition, these approximate solutions are “excellent approximations” (as mentioned in the statement of the main results) in a very strong sense: the distance to the exact solution dramatically decreases with a single iteration of Newton's method. In the context of eigenpair computations this concept is settled as follows.

Definition 2.11 Given $(A, \lambda, v) \in \mathcal{W}$ we say that $(\zeta, w) \in \mathbb{C} \times (\mathbb{C}^n \setminus \{0\})$ is an *approximate eigenpair* of A with associated eigenpair (λ, v) when for all $k \geq 1$ the k th iterate $N_A^k(\zeta, w)$ of the Newton map at (ζ, w) is well defined and satisfies

$$\text{dist}_A((N_A^k(\zeta, w)), (\lambda, v)) \leq \left(\frac{1}{2}\right)^{2^k-1} \text{dist}_A((\zeta, w), (\lambda, v)).$$

The following result estimates, in terms of the condition of an eigenpair, the radius of a ball of approximate eigenpairs associated to it. For a complete proof see [3, Theorem 5].

Theorem 2.12 *There is a universal constant $c_0 > 1/5$ with the following property. Let $(A, \lambda, v) \in \mathcal{W}$ with $\|A\|_F = 1$ and let $(\zeta, w) \in \mathbb{C} \times (\mathbb{C}^n \setminus \{0\})$. If*

$$\text{dist}_A((\lambda, v), (\zeta, w)) \leq \frac{c_0}{\mu(A, \lambda, v)},$$

then (ζ, w) is an approximate eigenpair of A with associated eigenpair (λ, v) .

It is a simple exercise to check that for any nonzero $z \in \mathbb{C}$, (ζ, w) is an approximate zero of A with associated zero (λ, v) if and only if $(z\zeta, w)$ is an approximate

zero of zA with associated zero $(z\lambda, v)$. So from the point of view of analyzing the effect of the Newton methods we may pick whatever scaling is convenient. For us it will be convenient to assume that $\|A\|_F = 1$ which we will do in the following.

PROOF OF THEOREM 2.12. Note that [3, Theorem 5] is the same result with $c_0 = 0.2881$, but the definition of the condition number in [3] is slightly different from ours. More exactly, if we denote by $\kappa(A, \lambda, v)$ the condition number defined in [3] then we have $\kappa(A, \lambda, v) = \max(1, \mu(A, \lambda, v))$. Theorem 2.12 is hence true with κ in the place of μ and $c_0 = 0.2881$. However, from Lemma 2.6 we know that $\mu(A, \lambda, v) \geq 2^{-1/2}$ which readily implies $\kappa(A, \lambda, v) \leq \sqrt{2}\mu(A, \lambda, v)$. Theorem 2.12 now follows from the fact that $0.2881 > \sqrt{2}/5$. \square

Remark 2.13 We note that $N_A(\zeta, w)$ can be computed from the matrix A and the pair (ζ, w) in $\mathcal{O}(n^3)$ operations, since the cost of this computation is dominated by that of inverting a matrix (or simply solving a linear system).

2.6 Gaussian Measures on $\mathbb{C}^{n \times n}$

Let $\sigma > 0$. We say that the complex random variable $Z = X + \sqrt{-1}Y$ has distribution $\mathcal{N}_{\mathbb{C}}(0, \sigma^2)$ when the real part X and the imaginary part Y are independent and identically distributed (i.i.d.) drawn from $\mathcal{N}(0, \frac{\sigma^2}{2})$, i.e., they are Gaussian centered random variables with variance $\frac{\sigma^2}{2}$.

If $Z \sim \mathcal{N}_{\mathbb{C}}(0, \sigma^2)$ then its density $\varphi: \mathbb{C} \rightarrow \mathbb{R}$ with respect to the Lebesgue measure is given by

$$\varphi(z) := \frac{1}{\pi\sigma^2} e^{-\frac{|z|^2}{\sigma^2}}.$$

We will write $v \sim \mathcal{N}_{\mathbb{C}^n}(0, \sigma^2)$ to indicate that the vector $v \in \mathbb{C}^n$ is random with i.i.d. coordinates drawn from $\mathcal{N}_{\mathbb{C}}(0, \sigma^2)$. Also, we say that $A \in \mathbb{C}^{m \times n}$ is (isotropic) Gaussian and we write $A \sim \mathcal{N}_{\mathbb{C}^{m \times n}}(0, \sigma^2)$, if its entries are i.i.d. Gaussian random variables. The resulting probability space is sometimes called the *Ginibre ensemble*.

If $\hat{A} \in \mathbb{C}^{m \times n}$ and $G \sim \mathcal{N}_{\mathbb{C}^{m \times n}}(0, \sigma^2)$, we say that the random matrix $A = G + \hat{A}$ has the *Gaussian distribution centered at \hat{A}* , and we write $A \sim \mathcal{N}_{\mathbb{C}^{m \times n}}(\hat{A}, \sigma^2)$. The density of this distribution is given by

$$\varphi_{m \times n}^{\hat{A}, \sigma}(A) := \frac{1}{(\pi\sigma^2)^{mn}} e^{-\frac{\|A - \hat{A}\|_F^2}{\sigma^2}}.$$

For conciseness, we will sometimes write $A \sim \mathcal{N}_{\mathbb{C}^{m \times n}}$ in the particular case where $\hat{A} = 0$ and $\sigma = 1$.

Crucial in our development is the following result giving a bound on the average condition for Gaussian matrices arbitrarily centered. Its statement is similar to the main technical result in [19, Thm. 3.6]. We will prove it in §7.

For technical reasons we will be interested in the following variation of μ :

$$\mu_F(A, \lambda, v) := \|A\|_F \|A_{\lambda, v}^{-1}\|_F$$

(note, we only replaced $\|A_{\lambda, v}^{-1}\|$ by $\|A_{\lambda, v}^{-1}\|_F$) and the corresponding

$$\mu_{F, \text{av}}(A) := \left(\frac{1}{n} \sum_{j=1}^n \mu_F^2(A, \lambda_j, v_j) \right)^{\frac{1}{2}}.$$

Theorem 2.14 For $\hat{A} \in \mathbb{C}^{n \times n}$ and $\sigma > 0$ we have

$$\mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}(\hat{A}, \sigma^2)} \left(\frac{\mu_{F, \text{av}}^2(A)}{\|A\|_F^2} \right) \leq \frac{n}{\sigma^2}.$$

Moreover, for A chosen with the uniform distribution $\mathcal{U}(\mathbb{S})$ in the unit sphere \mathbb{S} of $\mathbb{C}^{n \times n}$ we have:

$$\mathbb{E}_{A \sim \mathcal{U}(\mathbb{S})} (\mu_{F, \text{av}}^2(A)) \leq n^3.$$

Remark 2.15 (i) We note that no bound on the norm of \hat{A} is required in the first claim of Theorem 2.14. Indeed, using $\mu_{F, \text{av}}(sA) = \mu_{F, \text{av}}(A)$, it is easy to see that the assertion for a pair (\hat{A}, σ) implies the assertion for $(s\hat{A}, s\sigma)$, for any $s > 0$.

(ii) It is remarkable that if we change Gaussian matrices to some classes of structured matrices, the expected value of the condition number can be very high, see for example [10] and references therein.

2.7 Truncated Gaussians and smoothed analysis

For $T, \sigma > 0$, we define the *truncated Gaussian* $\mathcal{N}_{\mathbb{C}^{n \times n}, T}(0, \sigma^2)$ on $\mathbb{C}^{n \times n}$ to be the distribution given by the density

$$\rho_T^\sigma(A) = \begin{cases} \frac{\varphi_{n \times n}^{0, \sigma}(A)}{P_{T, \sigma}} & \text{if } \|A\|_F \leq T, \\ 0 & \text{otherwise,} \end{cases} \quad (13)$$

where $P_{T, \sigma} := \text{Prob}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}(0, \sigma^2)} \{\|A\|_F \leq T\}$, and, we recall, $\varphi_{n \times n}^{0, \sigma}$ is the density of $\mathcal{N}_{\mathbb{C}^{n \times n}}(0, \sigma^2)$. For the rest of this paper we fix the threshold $T := \sqrt{2}n$. The fact that $\|A\|_F^2$ is chi-square distributed with $2n^2$ degrees of freedom, along with [20, Corollary 6] yields the following result.

Lemma 2.16 We have $P_{T, \sigma} \geq \frac{1}{2}$ for all $0 < \sigma \leq 1$. \square

The space $\mathbb{C}^{n \times n}$ of matrices with the Frobenius norm and the space \mathbb{C}^{n^2} with the canonical Hermitian product are isomorphic as Hermitian product spaces. Hence, the Gaussian $\mathcal{N}_{\mathbb{C}^{n \times n}}(0, \sigma^2)$ on the former corresponds to the Gaussian $\mathcal{N}_{\mathbb{C}^{n^2}}(0, \sigma^2)$ on the latter, and we can deduce invariance of $\mathcal{N}_{\mathbb{C}^{n \times n}}(0, \sigma^2)$ under the action of \mathcal{U}_{n^2} (in addition to that for conjugation under \mathcal{U}_n discussed in §2.3), and the same is true for the truncated Gaussian. In particular, the pushforward of both distributions for the projection $\mathbb{C}^{n \times n} \setminus \{0\} \rightarrow \mathbb{S}$, $A \mapsto \frac{A}{\|A\|_F}$, is the uniform distribution $\mathcal{U}(\mathbb{S})$ (see [19, Chapter 2] for details) and we have

$$\mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}(0, \sigma^2)} F(A) = \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}, T}(0, \sigma^2)} F(A) = \mathbb{E}_{A \sim \mathcal{U}(\mathbb{S})} F(A), \quad (14)$$

for any measurable scale invariant function $F : \mathbb{C}^{n \times n} \rightarrow [0, \infty)$.

Complexity analysis has traditionally been carried out either in the *worst-case* or in an *average-case*. More generally, for a function $F : \mathbb{R}^m \rightarrow \mathbb{R}_+$ (some measure for the computational cost of solving an instance in \mathbb{R}^m), the former amounts to the evaluation of $\sup_{a \in \mathbb{R}^m} F(a)$ and the latter to that of $\mathbb{E}_{a \sim \mathcal{D}} F(a)$ for some probability distribution \mathcal{D} on \mathbb{R}^m . Usually, \mathcal{D} is taken to be the standard Gaussian in the input space. With the beginning of the century, Daniel Spielman and Shang-Hua Teng introduced a third form of analysis, *smoothed analysis* (see [53, 55] or [19, §2.2.7]), which is meant to interpolate between worst-case and average-case.

The idea is to replace the two operators above (supremum and expectation) by a combination of the two, namely,

$$\sup_{\hat{a} \in \mathbb{R}^m} \mathbb{E}_{a \sim \mathcal{D}(\hat{a}, \sigma)} F(a)$$

where $\mathcal{D}(\hat{a}, \sigma)$ is a distribution “centered” at \hat{a} having σ as a measure of dispersion. A typical example is the Gaussian $\mathcal{N}(\hat{a}, \sigma^2)$. Another example, used for scale invariant functions F , is the uniform measure on a spherical cap centered at \hat{a} and with angular radius σ on the unit sphere $\mathbb{S}(\mathbb{R}^m)$ (reference [19] exhibits smoothed analyses for both examples of distribution). In this paper we will perform a smoothed analysis with respect to a truncated Gaussian. More precisely, we will be interested in quantities of the form

$$\sup_{\hat{A} \in \mathbb{C}^{n \times n}} \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}, T}(\hat{A}, \sigma^2)} F(A)$$

where F will be a measure of computational cost for the eigenpair problem. We note that, in addition to the usual dependence on n , this quantity depends also on σ and tends to ∞ when σ tends to 0. When F is scale invariant, as in the case of μ_{av} or μ_{max} , it is customary to restrict attention to matrices of norm 1. That is, to study the following quantity:

$$\sup_{\hat{A} \in \mathbb{S}} \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}, T}(\hat{A}, \sigma^2)} F(A). \quad (15)$$

2.8 The eigenpair continuation algorithm

We are ready to describe the main algorithmic construct in this paper. When dealing with algorithms it will be more convenient to view the solution variety as the corresponding subset of $\mathbb{C}^{n \times n} \times \mathbb{C} \times (\mathbb{C}^n \setminus \{0\})$, which, abusing notation, we still denote by \mathcal{V} .

Given two matrices $B_0, B \in \mathbb{S}$, $B \neq \pm B_0$, let $\alpha := d_{\mathbb{S}}(B_0, B) \in (0, \pi)$ be the spherical distance (i.e. the angle) from B_0 to B , and let

$$\mathcal{L}_{B_0, B} = \{B_s : 0 \leq s \leq \alpha\} \quad (16)$$

be the portion of the great circle in \mathbb{S} , parametrized by arc-length, joining B_0 and B , so $B_\alpha = B$. By abuse of notation, for any $A_0, A \in \mathbb{C}^{n \times n}$ such that A_0, A are not \mathbb{R} -linearly dependent, we simply write

$$d_{\mathbb{S}}(A_0, A) := d_{\mathbb{S}}\left(\frac{A_0}{\|A_0\|_F}, \frac{A}{\|A\|_F}\right) \text{ and } \mathcal{L}_{A_0, A} := \mathcal{L}_{\frac{A_0}{\|A_0\|_F}, \frac{A}{\|A\|_F}}.$$

The following definition plays a distinguished role in the continuation algorithm. It uses a constant c_* which we will later take to be 10^{-4} (cf. §5.1).

Definition 2.17 If $(A, \lambda, v) \in \mathcal{W}$, $(\zeta, w) \in \mathbb{C} \times (\mathbb{C}^n \setminus \{0\})$ with $\text{dist}_A((\zeta, w), (\lambda, v)) \leq \frac{c_*}{\mu_{(A, \lambda, v)}}$, then we say that (ζ, w) is a *certified approximate eigenpair* of A (with associated eigenpair (λ, v)). The *initial neighborhood* of the set \mathcal{W} is the set

$$\widetilde{\mathcal{W}} := \left\{ (A, \zeta, w) \mid \exists (\lambda, v) \text{ s.t. } (A, \lambda, v) \in \mathcal{W} \text{ and } \text{dist}_A((\zeta, w), (\lambda, v)) \leq \frac{c_*}{\mu_{\max}(A)} \right\}.$$

The term *certified* in Definition 2.17 is justified by Theorem 2.12.

Suppose that we are given an *initial triple* $(B_0, \zeta_0, w_0) \in \widetilde{\mathcal{W}}$, $B_0 \in \mathbb{S}$ and an input matrix $B \in \mathbb{S} \setminus \{\pm B_0\}$. Let $(\lambda_0, v_0) \in \mathbb{C} \times \mathbb{P}(\mathbb{C}^n)$ be the exact eigenpair of B_0 associated to (ζ_0, v_0) . As a consequence of the inverse function theorem, if $(\mathcal{L}_{B_0, B} \setminus \{B_0\}) \cap \Sigma = \emptyset$, then the map $s \mapsto B_s$ can be uniquely extended to a continuous map

$$[0, \alpha] \rightarrow \mathcal{V}, \quad s \mapsto (B_s, \lambda_s, v_s), \quad (17)$$

We call this map the *lifting* of $\mathcal{L}_{B_0, B}$ with origin (B_0, λ_0, v_0) . We can try to approximate the eigenpair $(\lambda_\alpha, v_\alpha)$ of B by following the lifting of $\mathcal{L}_{B_0, B}$. To this end we can differentiate $B_s v_s - \lambda_s v_s = 0$ w.r.t. s . This produces an Initial Value Problem (IVP) whose solution can be approximated by any standard numerical ODE solver. The main ingredient for the complexity estimate is the number of points in the discretization of $[0, \alpha]$ needed to approximate the solution of the IVP.

Formalizing this idea to get an actual guarantee of convergence is a nontrivial task; only a non-constructive method has been described in [3] following the ideas in [44]. We now describe how to algorithmically construct a numerically stable

method for this task: subdivide the interval $[0, \alpha]$ into subintervals with extremities at $0 = s_0 < s_1 < \dots < s_K = \alpha$ and successively compute approximations (ζ_i, w_i) of (λ_{s_i}, v_{s_i}) , starting with (ζ_0, w_0) and then using Newton's method. To ensure that these are good approximations, we actually want to ensure that for all i , (ζ_i, w_i) is an approximate eigenpair of $B_{s_{i+1}}$. Figure 1 attempts to convey the general idea.

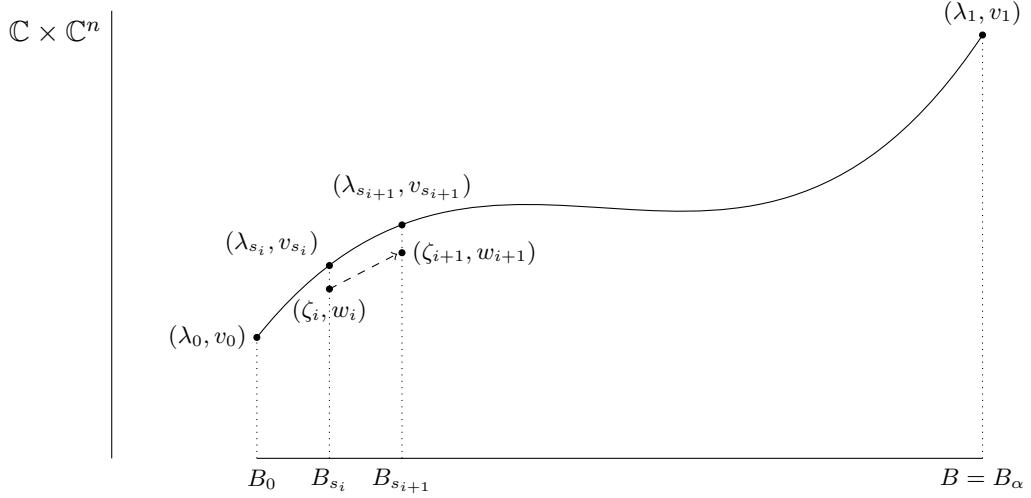


Figure 1: The continuation of the solution path.

The algorithmic counterpart of this idea is the following.

Algorithm 1 Path-follow

Input: $A, A_0 \in \mathbb{C}^{n \times n} \setminus \{0\}$, and $(\zeta_0, w_0) \in \mathbb{C} \times \mathbb{C}^n$

Preconditions: $\|w_0\| = 1$, $\mathbb{R}A_0 \neq \mathbb{R}A$, (ζ_0, w_0) is a certified approximate eigenpair of A_0 (with some associated eigenpair (λ_0, v_0)).

```

redefine  $A_0 := A_0/\|A_0\|_F$ ,  $A_1 := A/\|A\|_F$ 
 $\alpha := d_{\mathbb{S}}(A_0, A_1)$ ,  $s_0 := 0$ ,  $B := A_0$ ,  $i = 0$ 
repeat
     $\dot{A} :=$  unit tangent vector (in the direction of the
        parametrization) to  $\mathcal{L}_{A_0, A_1}$  at  $B$ 
     $\Delta s := \text{Choose\_step}(B, \dot{A}, \zeta_i, w_i)$ 
     $s_{i+1} := \min\{\alpha, s_i + \Delta s\}$ 
     $B :=$  point in  $\mathcal{L}_{A_0, A_1}$  with  $d_{\mathbb{S}}(A_0, B) = s$ 
     $(\zeta_{i+1}, w_{i+1}) := N_B^3(\zeta_i, w_i)$  (three Newton iterations)
    if  $|\zeta_{i+1}| > 1$  then  $\zeta_{i+1} = \zeta_{i+1}/|\zeta_{i+1}|$ .
     $i := i + 1$ 
until  $s = \alpha$ 
return  $(\zeta', w') = (\|A\|_F \zeta_i, w_i)$ 

```

Output: $(\zeta', w') \in \mathbb{C} \times \mathbb{C}^n$

Postconditions: The execution halts if the lifting (A_t, λ_t, v_t) , $0 \leq t \leq d_{\mathbb{S}}(A_0, A)$, of $\mathcal{L}_{A_0, A}$ at (λ_0, v_0) doesn't meet Σ' . In this case, (ζ', w') is a certified approximate eigenpair of A with associated eigenpair (λ_1, v_1) .

Remark 2.18 Note that in particular any triple $(A_0, \lambda_0, v_0) \in \mathcal{W}$ is an acceptable starting point for Path-follow for generic $A \in \mathbb{C}^{n \times n}$.

Algorithm Path-follow is unambiguous except for the subroutine Choose_step, which will be described at the end of §5.2 below. We next state the main results for it. Recall, the point $s_\ell \in [0, \alpha]$ is the value of s generated by Path-follow at the ℓ th iteration. For $K \in \mathbb{R}$ we denote by $\lceil K \rceil$ the smallest integer which is greater than or equal to K .

Theorem 2.19 Suppose that $\mathcal{L}_{A_0, A} = \{B_s\}_{s \in [0, \alpha]}$ (so $B_0 = A_0/\|A_0\|_F$ and $B_\alpha = A/\|A\|_F$) and assume that $\mathcal{L}_{A_0, A} \cap \Sigma = \emptyset$. Then the algorithm Path-follow stops after at most $\lceil K \rceil$ steps where K is given by

$$K := K(A, A_0, \zeta_0, w_0) := C \int_0^\alpha \mu(B_s, \lambda_s, v_s) \|(\dot{B}_s, \dot{\lambda}_s, \dot{v}_s)\| ds.$$

Here the pairs (λ_s, v_s) are given by (17).

More generally, let $q \in \mathbb{Z}, q \geq 0$. Then algorithm **Path-follow** stops after at most (the smallest integer greater than or equal to)

$$q + C \int_{s_q}^{\alpha} \mu(B_s, \lambda_s, v_s) \|(\dot{B}_s, \dot{\lambda}_s, \dot{v}_s)\| ds$$

steps. The returned pair (ζ, w) is an approximate eigenpair of A with associated eigenpair $(\lambda_\alpha, v_\alpha)$. Here, $C \leq 3000$ is a universal constant.

For $0 \leq a < b \leq \alpha$, the quantity

$$L_{\mu,a,b}(B_s, \lambda_s, v_s) = \int_a^b \mu(B_s, \lambda_s, v_s) \|(\dot{B}_s, \dot{\lambda}_s, \dot{v}_s)\| ds \quad (18)$$

in Theorem 2.19 is the length of the curve $\{(B_s, \lambda_s, v_s) : a \leq s \leq b\}$ in the so-called *condition metric*. This is the metric that is obtained by pointwise multiplying the natural metric in $\mathbb{S} \times \mathbb{C} \times \mathbb{P}(\mathbb{C}^n)$ by the condition number squared. We call $L_{\mu,a,b}(B_s, \lambda_s, v_s)$ the *condition length* of this curve.

The proof of Theorem 2.19 is given in §5.3.

Remark 2.20 From Proposition 2.7, we have

$$L_{\mu,s_q,\alpha}(B_s, \lambda_s, v_s) \leq \sqrt{6} \int_{s_q}^{\alpha} \mu^2(B_s, \lambda_s, v_s) ds.$$

The following result gives an alternative bound to the number of steps.

Corollary 2.21 Let $A_0, A \in \mathbb{C}^{n \times n}$ be \mathbb{R} -linearly independent and consider the path $A_t = (1-t)A_0 + tA$ which satisfies $A_1 = A$. If $\mathcal{L}_{A_0, A} \cap \Sigma = \emptyset$ then, for any $q \in \mathbb{Z}, q \geq 0$, the algorithm **Path-follow** stops after at most $\lceil K \rceil$ steps where

$$K := q + \sqrt{6}C \|A_0\|_F \|A_1\|_F \int_{t_q}^1 \frac{\mu^2(A_t, \lambda_t, v_t)}{\|A_t\|_F^2} dt$$

steps, with

$$t_q := \frac{\|A_0\|_F}{\|A_1\|_F (\sin \alpha \cot s_q - \cos \alpha) + \|A_0\|_F}.$$

Here, λ_t, v_t are the eigenvalue and eigenvector of A_t defined by continuation.

PROOF. From Theorem 2.19 and Remark 2.20, letting $C' = \sqrt{6}C$, the number of steps is at most the smallest integer greater than or equal to

$$q + C' \int_{s_q}^{\alpha} \mu^2(B_s, \lambda_{B_s}, v_{B_s}) ds = q + C' \int_{s_q}^{\alpha} \mu^2(B_s, \lambda_{B_s}, v_{B_s}) \|\dot{B}_s\|_F ds,$$

where λ_{B_s} and v_{B_s} are the eigenvalue and eigenvector of B_s defined by continuation. Now, reparametrizing that spherical segment by $C_t = A_t/\|A_t\|_F$, $t \in [0, 1]$, the integral does not change. The quantity above is thus equal to

$$q + C' \int_{t_q}^1 \mu^2(C_t, \lambda_{C_t}, v_{C_t}) \|\dot{C}_t\|_F dt \underset{\text{Rmk. 2.5}}{=} q + C' \int_{t_q}^1 \mu^2(A_t, \lambda_t, v_t) \left\| \frac{d}{dt} \left(\frac{A_t}{\|A_t\|_F} \right) \right\| dt.$$

Substituting $\dot{A}_t = A_1 - A_0$ and $A_t = (1-t)A_0 + tA_1$ in this last formula and with some elementary computations (see [19, Lemma 17.5]) we conclude that

$$\left\| \frac{d}{dt} \left(\frac{A_t}{\|A_t\|_F} \right) \right\| \leq \frac{\|A_0\|_F \|A_1\|_F}{\|A_t\|_F^2}.$$

The corollary follows (the value of t_q is obtained by the reparametrization from s_q). \square

The inequality of Remark 2.20 implies that (up to constants) the upper bound for the number of steps by an algorithm in terms of the condition length as in Theorem 2.19 is smaller than the upper bound in terms of the integral of the squared condition number as in Corollary 2.21. A similar situation applies in the context of polynomial system root finding. In this case implementations exist in both contexts see [49, 18, 11, 23]. The condition length algorithm is more subtle and the proof of correctness more difficult, both for the polynomial system and eigenpair cases. So the temptation is to present condition number squared algorithms. Here for the first time we present a quantitative estimate of the improvement which is significant and which justifies presenting the more complex condition length algorithm (another estimate of this kind was established shortly after in [4]). In Theorem 2.29 a randomized algorithm is studied. The upper bound given by the condition length algorithm is $O(n^2)$ while an algorithm with complexity given by the condition number squared would give $O(n^3)$.

In our main results we are interested in the cost of our algorithms over random matrices A . The following quantity —the expected number of iterations of Path-follow for a given initial triple $(A_0, \lambda_0, v_0) \in \mathcal{W}$ — becomes essential,

$$\text{Avg_Num_Iter}(A_0, \lambda_0, v_0) := \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}} K(A, A_0, \lambda_0, v_0).$$

We can also consider the smoothed number of iterations of Path-follow that results by drawing instead the input matrix A from $\mathcal{N}_T(\hat{A}, \sigma^2)$ where $\hat{A} \in \mathbb{S}$ is arbitrary. We thus define

$$\text{Smd_Num_Iter}(A_0, \lambda_0, v_0, \sigma) := \sup_{\hat{A} \in \mathbb{S}} \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}, T}(\hat{A}, \sigma^2)} K(A, A_0, \lambda_0, v_0).$$

Proposition 2.22 *For $(A_0, \lambda_0, v_0) \in \mathcal{W}$ we have*

$$\text{Avg_Num_Iter}(A_0, \lambda_0, v_0) = \mathcal{O}(n^4 \mu^2(A_0, \lambda_0, v_0))$$

and, for all $\sigma \in (0, 1]$,

$$\text{Smd_Num_Iter}(A_0, \lambda_0, v_0, \sigma) = \mathcal{O}\left(\frac{n^4 \mu^2(A_0, \lambda_0, v_0)}{\sigma^2}\right).$$

Remark 2.23 Note that Proposition 2.22 ensures that Path-follow halts in finite time with probability 1 even if $A_0 \in \Sigma$, as long as $\mu(A_0, \lambda_0, v_0)$ is finite. The algorithm does a first step that depends on $\mu(A_0, \lambda_0, v_0)$ only and advances to B_{s_1} with $s_1 > 0$. After that, the fact that the real codimension of Σ in $\mathbb{C}^{n \times n}$ is 2 (shown in Proposition 2.3) ensures that, almost surely, $(\mathcal{L}_{A_0, A} \setminus \{A_0\}) \cap \Sigma = \emptyset$. Therefore, with probability 1, none of the matrices B_s is in Σ and the integral over $[s_1, \alpha]$ in Theorem 2.19 is finite. Also, note that one needs not to have the exact eigenpair (λ_0, v_0) , but it suffices to have a certified approximate eigenpair of A_0 with associated eigenpair (λ_0, v_0) . In particular, it suffices to have $(A_0, \lambda_0, w_0) \in \widetilde{\mathcal{W}}$.

To compute all the eigenpairs from an initial matrix $A_0 \in \mathbb{C}^{n \times n} \setminus \Sigma$ and its eigenpairs $(\lambda^{(1)}, v^{(1)}), \dots, (\lambda^{(n)}, v^{(n)})$ we may proceed by following the n paths corresponding to taking these eigenpairs in the initial triples. In this case, we will be interested in the quantities

$$\begin{aligned} \text{Avg_Num_Iter_All}(A_0) &:= \sum_{i=1}^n \text{Avg_Num_Iter}(A_0, \lambda^{(i)}, v^{(i)}) \\ &= \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}} \sum_{i=1}^n K(A, A_0, \lambda^{(i)}, v^{(i)}), \end{aligned}$$

and

$$\text{Smd_Num_Iter_All}(A_0, \sigma) := \sup_{\hat{A} \in \mathbb{S}} \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}, T}(\hat{A}, \sigma^2)} \sum_{i=1}^n K(A, A_0, \lambda^{(i)}, v^{(i)}).$$

For these quantities we prove the following result.

Proposition 2.24 For $A_0 \in \mathbb{C}^{n \times n} \setminus \Sigma$ we have

$$\text{Avg_Num_Iter_All}(A_0) = \mathcal{O}(n^4 \mu_{\max}^2(A_0))$$

and, for all $\sigma \in (0, 1]$,

$$\text{Smd_Num_Iter_All}(A_0, \sigma) = \mathcal{O}\left(\frac{n^4 \mu_{\max}^2(A_0)}{\sigma^2}\right).$$

We prove Propositions 2.22 and 2.24 in Section 8. Note that the latter does not immediately follow from the former since in general

$$\sum_{i=1}^n \mu^2(A_0, \lambda^{(i)}, v^{(i)}) >> \mu_{\max}^2(A_0).$$

2.9 Initial triples and global algorithms

The Path-follow routine assumes an initial triple (A_0, λ_0, v_0) at hand. We next deal with this issue. We first consider the case of computing a single eigenpair. In this case we consider the diagonal matrix H whose diagonal entries are $(1, 0, \dots, 0)$ and its well-posed eigenpair $(1, e_1)$.

Algorithm 2 Single_Eigenpair

Input: $A \in \mathbb{C}^{n \times n}$

$(\zeta, w) := \text{Path-follow}(A, H, 1, e_1)$

Output: $(\zeta, w) \in \mathbb{C} \times \mathbb{C}^n$

Postconditions: The execution halts if the lifting of $\mathcal{L}_{H,A}$ at $(1, e_1)$ doesn't meet Σ' . In this case, the returned (ζ, w) is a certified approximate eigenpair of A .

We can formally state (and prove) the first of our main results. To this end, we define the average cost $\text{Avg_Cost}(n)$ of Single_Eigenpair to be the average (over the input matrix A) of the number of arithmetic operations performed by the algorithm. We similarly define its smoothed cost $\text{Smd_Cost}(n, \sigma)$.

Theorem 2.25 *Algorithm Single_Eigenpair returns (almost surely) an approximate eigenpair of its input $A \in \mathbb{C}^{n \times n}$. Its average cost satisfies*

$$\text{Avg_Cost}(n) = \mathcal{O}(n^7).$$

For every $0 < \sigma \leq 1$, its smoothed cost satisfies

$$\text{Smd_Cost}(n, \sigma) = \mathcal{O}\left(\frac{n^7}{\sigma^2}\right).$$

PROOF. Lemma 2.10 and the fact that $\|H\|_F = 1$ imply that $\mu(H, 1, e_1) = 1$. The statement is then a consequence of Proposition 2.22 and the fact that the average cost is obtained by multiplying $\text{Avg_Num_Iter}(H, 1, e_1)$ by the cost $\mathcal{O}(n^3)$ of each iteration. \square

Remark 2.26 The triple $(H, 1, e_1)$ is the version, in our context, of the initial pair proposed by Shub and Smale in [49] for the computation of zeros of polynomial systems. In this later context, the problem of showing that one can efficiently follow linear homotopies with this initial pair remains open.

The fact that any other eigenpair of H is ill-posed prevents us from using them to compute other eigenpairs of A . If we want to compute all the eigenpairs of A we will need to consider a different approach.

To do so, for any fixed $n \geq 2$, let

$$D = \text{diag}(\eta_1, \dots, \eta_n), \quad (19)$$

where the η_i are points in the unit side hexagonal lattice chosen in such an order that $0 = |\eta_1| \leq \dots \leq |\eta_n|$. We recall, the hexagonal lattice is the set of points of the form

$$Q \begin{pmatrix} a \\ b \end{pmatrix}, \quad \text{where} \quad Q = \begin{pmatrix} 1 & 1/2 \\ 0 & \sqrt{3}/2 \end{pmatrix}, \quad a, b \in \mathbb{Z}.$$

Lemma 2.27 *We have $\mu_{\max}(D) \leq \frac{\sqrt{3/2}}{\pi} n + o(n)$.*

PROOF. We first find a real number $r > 0$ with the property that the disk $\mathbb{D}(r)$ of radius r contains at least n points in the hexagonal lattice. To do so, note that a lattice point $Q \begin{pmatrix} a \\ b \end{pmatrix}$ is in $\mathbb{D}(r)$ if and only if

$$\begin{pmatrix} a \\ b \end{pmatrix} \in Q^{-1}\mathbb{D}(r).$$

Now, the singular values of Q^{-1} are $\sqrt{2}$ and $\sqrt{2/3}$, so $Q^{-1}\mathbb{D}(r)$ is an ellipse of area $2\pi r^2/\sqrt{3}$ and maximal radius $\sqrt{2}r$. Dividing by the smallest integer N that is greater than $2\sqrt{2}r$ and translating the resulting ellipse to have center $(1/2, 1/2)$, we look for points of the form $(a/N, b/N)$ with $a, b \in \{0, \dots, N-1\}$ which lie inside an ellipse of area

$$\frac{2\pi r^2}{\sqrt{3}N^2}$$

contained in $[0, 1]^2$. This is a particular instance of the problem of counting lattice points in semialgebraic sets, a well studied problem for which a quite complete solution is for example [14, Th. 3, p. 327]. We conclude that the number of points in the hexagonal lattice in $\mathbb{D}(r)$ is at least

$$\frac{2\pi r^2}{\sqrt{3}} - 2N \geq \frac{2\pi r^2}{\sqrt{3}} - 4\sqrt{2}r - 2.$$

In particular, we can find n lattice points in a disk of radius $r = 3^{1/4}/(2\pi)^{1/2}n^{1/2} + o(n^{1/2})$. Moreover, around each lattice point we can place a circle of radius $1/2$ without overlappings. From the mean value equality for analytic functions, for any $z \in \mathbb{C}$ and $\varepsilon > 0$ we have,

$$|z|^2 = |z^2| = \left| \frac{1}{\pi\varepsilon^2} \int_{|y-z|<\varepsilon} y^2 dy \right| \leq \frac{1}{\pi\varepsilon^2} \int_{|y-z|<\varepsilon} |y|^2 dy.$$

We thus have

$$\|D\|_F^2 = \sum_{j=1}^n |\eta_j|^2 \leq \sum_z |z|^2 \leq \frac{4}{\pi} \sum_z \int_{|y-z|<1/2} |y|^2 dy \leq \frac{4}{\pi} \int_{|y|<r+1/2} |y|^2 dy,$$

where z runs over all lattice points contained in $\mathbb{D}(r)$. Solving the last integral yields

$$\|D\|_F^2 \leq 2r^4 + o(r^4) = \frac{3n^2}{2\pi^2} + o(n^2).$$

Finally, from Lemma 2.10 we conclude that

$$\mu_{\max}^2(D) = \|D\|_F^2 \leq \frac{3n^2}{2\pi^2} + o(n^2). \quad \square$$

We now put together the continuation algorithm `Path-follow` and this specific initial triple.

Algorithm 3 All_Eigenpairs

Input: $A \in \mathbb{C}^{n \times n}$

compute D and η_j as defined in (19)
for $j \in \{1, \dots, n\}$ **do**
 $(\zeta_j, w_j) := \text{Path-follow}(A, D, \eta_j, e_j)$

Output: $((\zeta_1, w_1), \dots, (\zeta_n, w_n)) \in (\mathbb{C} \times \mathbb{C}^n)^n$

Postconditions: The algorithm halts if $\mathcal{L}_{D,A} \cap \Sigma = \emptyset$. In this case, the pairs (ζ_j, w_j) are certified approximate eigenpairs of A with pairwise different associated eigenpairs.

We can now state (and prove) the second of our main results.

Theorem 2.28 *Algorithm All_Eigenpairs returns (almost surely) n approximate eigenpairs of its input $A \in \mathbb{C}^{n \times n}$, with pairwise different associate eigenpairs. Its average cost satisfies*

$$\text{Avg_Cost}(n) = \mathcal{O}(n^9).$$

For every $\sigma \leq 1$ its smoothed cost satisfies

$$\text{Smd_Cost}(n, \sigma) = \mathcal{O}\left(\frac{n^9}{\sigma^2}\right).$$

PROOF. It easily follows from Lemma 2.27 and Proposition 2.24. As in the proof of Theorem 2.25, we recall that the cost of each iteration is $O(n^3)$. \square

2.10 Computer experiments

Algorithm `Single_Eigenpair` was implemented in Matlab by Liu Yiting, and this implementation was used to get an empirical estimate of its average number of iterations and average cost.

To estimate these averages she generated 200 random complex matrices $A_j \in \mathbb{C}^{n \times n}$ for $n = 4, 8, 16, 32$. For each n and each j she computed an eigenpair for A_j and recorded the number I_{nj} of iterations performed by `Single_Eigenpair` to do so. Then she computed the (empirical) average number of iterations $\text{Iter}(n) = \frac{1}{200} \sum_{j=1}^{200} I_{nj}$. She did the same for $n = 64$ but only with 30 matrices. She obtained, respectively, 1571.4, 3464.9, 6410.4, 9390.6, and 13941.

In order to estimate the growth of $\text{Iter}(n)$ with n she did a linear regression between $\log_2 \text{Iter}(n)$ and $\log_2 n$ for $n = 4, 8, 16, 32$ and 64. The regression line obtained was $\log_2 \text{Iter}(n) \approx 0.7737(\log_2 n) + 9.3026$. The value of the regression coefficient, close to 1, shows that this line is a very good fit to the functional dependence $\log_2 \text{Iter}(n) = f(\log_2 n)$. Taking exponential with base 2 on both sides, we get $\text{Iter}(n) \approx 632 n^{0.78}$, and therefore, an empirical estimate of $\mathcal{O}(n^{3.78})$ for $\text{Avg_Cost}(n)$, substantially better than the $\mathcal{O}(n^7)$ in Theorem 2.25 (but still worse than the empirical behavior of the eigensolvers currently used).

2.11 Randomized algorithms

In this section we follow the ideas in [13] adapting them to the case of eigenpair computations. Consider the probability distribution \mathcal{D} in the solution variety \mathcal{V} defined via the following procedure:

randomly choose $A_0 \sim \mathcal{N}_{\mathbb{C}^{n \times n}}$ (20)
 randomly choose one eigenpair (λ_0, v_0) of A_0

Next assume that we have a routine `draw_from_D` to draw triples (A_0, λ_0, v_0) from the distribution \mathcal{D} on \mathcal{V} . Then we can consider the following algorithm.

Algorithm 4 `Random_initial_triple` (scheme)

Input: $A \in \mathbb{C}^{n \times n}$

$(A_0, \lambda_0, v_0) := \text{draw_from_}\mathcal{D}$
 $(\zeta, w) := \text{Path-follow}(A, A_0, \lambda_0, v_0)$

Output: $(\zeta, w) \in (\mathbb{C} \times \mathbb{C}^n)^n$

Postconditions: The execution halts if the lifting of $\mathcal{L}_{A_0, A}$ at (λ_0, v_0) doesn't meet Σ' . In this case, (ζ, w) is an approximate eigenpair of A .

The interest of this algorithmic scheme is that we can prove better bounds (than those in Theorem 2.25) for the number of iterations of **Path-follow**.

Theorem 2.29 *The expected average number of homotopy steps of Path-follow in Algorithm 4 satisfies*

$$\mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}} \mathbb{E}_{(A_0, \lambda_0, v_0) \sim \mathcal{D}} K(A, A_0, \lambda_0, v_0) \leq 4Cn^2,$$

where C is as in Theorem 2.19.

Theorem 2.29, which will be proved in Section 9, brings to focus the need of an implementation of `draw_from_D`. It must be noted though that a direct implementation is not possible since the second line in (20) (choosing (λ_0, v_0) at random) implicitly requires solving an EVP problem, the very question that this article is attempting to solve! This is a similar situation to that solved in [12, 13], where a random polynomial system and one of its zeros at random had to be chosen. It is also similar to that dealt with in [5] for the computation of eigenpairs of Hermitian matrices. A version of the proof of Theorem 2.29 with the Gaussian Unitary Ensemble replacing the Gaussian distribution, actually yields an improvement over the main result in [5]. (We note in passing that the case of Hermitian matrices is one of the few instances where a proof of convergence of a practical eigensolver, such as the Francis-Kublanovskaya's QR, is actually known [7].)

It must also be noted that following the method in [34] for the polynomial system case Algorithm 4 can probably be de-randomized without increasing the complexity.

Corollary 2.30 *In the case of Hermitian matrices, the expected average number of homotopy steps of Path-follow in Algorithm 4 (with the randomization algorithm in [5] in place of `draw_from_D`) is $\mathcal{O}(n^2)$. This yields an expected cost of $\mathcal{O}(n^5)$ for the computation of one eigenpair and of $\mathcal{O}(n^6)$ for the computation of all of them. Here the input matrix H is drawn from the Gaussian Unitary Ensemble. \square*

Following the ideas in those papers, we note that Theorem 2.29 would yield an algorithm with average running time $\mathcal{O}(n^5)$ if we could find some collection of probability spaces Ω_n and functions $\psi_n : \Omega_n \rightarrow \mathcal{V}_n$, $n \geq 2$, such that:

1. Choosing $\omega \in \Omega_n$ can be done starting with a number of random choices of numbers with the $\mathcal{N}_{\mathbb{C}}$ distribution, and performing some arithmetic operations on the results, the total expected running time being at most $\mathcal{O}(n^5)$.
2. Given $\omega \in \Omega_n$, $\psi_n(\omega)$ is computable in average time $\mathcal{O}(n^5)$, that is, the expected number of arithmetic operations for computing $\psi_n(\omega)$ must be $\mathcal{O}(n^5)$.
3. Choosing ω at random in Ω_n and computing $(A_0, \lambda_0, v_0) = \psi_n(\omega)$ is equivalent to choosing $A_0 \sim \mathcal{N}_{\mathbb{C}^{n \times n}}$ at random and choosing at random (λ_0, v_0) such that

$Av_0 = \lambda_0 v_0$. That is, for any measurable mapping $\phi : \mathcal{V}_n \rightarrow [0, \infty]$ we must have

$$\mathbb{E}_{\omega \sim \Omega_n} (\phi(\psi_n(\omega))) = \mathbb{E}_{A_0 \sim \mathcal{N}_{\mathbb{C}^{n \times n}}} \left(\frac{1}{n} \sum_{\lambda_0, v_0 : A_0 v_0 = \lambda_0 v_0} \phi(A_0, \lambda_0, v_0) \right), \quad (21)$$

so that we can apply this equality to

$$\phi(A_0, \lambda_0, v_0) = \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}} (K(A, A_0, \lambda_0, v_0))$$

and apply Theorem 2.29.

Unfortunately, we are not able to produce a collection of probability spaces Ω_n and functions ψ_n as described above. However, we will prove that relaxing (21) to the following less restrictive situation is actually possible: instead of demanding the equality in (21) we can just demand an inequality where the right-hand term is multiplied by some polynomial in n . Moreover, we do not need (21) to hold for every measurable function ϕ since all the interesting functions for the EVP problem are projective functions, invariant under the action of the unitary group. We can thus relax (21) to hold only with a polynomially bounded inequality, and for unitary invariant projective functions. Proving that this can actually be done is our goal now. (Meanwhile our ideas have been adapted to the more general setting of eigenpairs of polynomial systems in [15].)

We start by defining Ω_n and ψ_n . Consider the classical Stiefel manifold consisting of orthonormal $(n-1)$ -frames in \mathbb{C}^n , given by

$$\mathcal{S}_{n-1}(\mathbb{C}^n) = \{Q \in \mathbb{C}^{n \times (n-1)} : Q^* Q = I_{n-1}\},$$

endowed with its natural probability measure given by the restriction of the Frobenius Hermitian structure to the tangent bundle.

For every $n \geq 2$, let

$$\mathcal{A}_n := \{(M, Q) : \ker(M) = \ker(Q^*)\} \subseteq \mathbb{C}^{(n-1) \times n} \times \mathcal{S}_{n-1}(\mathbb{C}^n). \quad (22)$$

In other words, \mathcal{A}_n consists of pairs of matrices M, Q such that the columns of Q form an orthonormal basis of the complement of $\ker(M)$. The set \mathcal{A}_n has a natural probability measure $\mu_{\mathcal{A}_n}$ given by

$$\mu_{\mathcal{A}_n}(X) := \mathbb{E}_{M \sim \mathcal{N}_{\mathbb{C}^{(n-1) \times n}}} \left(\mathbb{E}_{Q : (M, Q) \in \mathcal{A}_n} (\mathbb{1}_X(M, Q) dQ) \right),$$

for measurable sets $X \subseteq \mathcal{A}_n$. Here, Q has the uniform distribution in the compact manifold $\{Q \in \mathcal{S}_{n-1}(\mathbb{C}^n) : (M, Q) \in \mathcal{A}_n\}$ and $\mathbb{1}_X$ is the indicator function of X .

Definition 2.31 Let

$$\Omega_n := \{(\lambda, w, (M, Q)) : 2\Re(\bar{\lambda}\text{tr}(MQ)) \leq 1 - |\lambda|^2(n-1)\} \subseteq \mathbb{C} \times \mathbb{C}^{n-1} \times \mathcal{A}_n$$

be endowed with the product measure μ_{Ω_n} , normalized to have total unit mass (see Remark 2.32 below). Then, let

$$\psi_n(\lambda, w, M, Q) = \left(\begin{pmatrix} \lambda & w^* \\ 0 & MQ + \lambda I_{n-1} \end{pmatrix}, \lambda, e_1 \right).$$

Remark 2.32 An explicit description of the product measure μ_{Ω_n} is as follows:

$$\mu_{\Omega_n}(Y) := C_n \mathbb{E}_{\lambda, w} (\mu_{\mathcal{A}_n}(\{(M, Q) : (\lambda, w, M, Q) \in Y\})),$$

for measurable sets $Y \subseteq \Omega_n$, where $\lambda \sim \mathcal{N}_{\mathbb{C}}$, $w \sim \mathcal{N}_{\mathbb{C}^{n \times 1}}$ and C_n is a normalizing constant given by

$$C_n^{-1} = \mathbb{P}_{(\lambda, M, Q) \in \mathbb{C} \times \mathcal{A}_n} (2\Re(\bar{\lambda}\text{tr}(MQ)) \leq 1 - |\lambda|^2(n-1)). \quad (23)$$

Our last main result (see Section 10 for a proof) is the following.

Theorem 2.33 Let Ω_n and ψ_n be as in Definition 2.31 for all $n \geq 2$. Then:

1. Choosing $\omega \in \Omega_n$ can be done by choosing $2n^2 - 2n + 1$ numbers with the $\mathcal{N}_{\mathbb{C}}$ distribution, checking a test which involves the computation of a Moore-Penrose inverse and computing two QR decompositions. This process may be repeated as a function of the test's outcome, but the expectation of the number of times the test is performed is at most C_n . The total expected running time is $\mathcal{O}(n^3 C_n)$.
2. Given $\omega \in \Omega_n$, computing $\psi_n(\omega)$ can be done with running time $\mathcal{O}(n^3)$.
3. For any unitarily invariant measurable mapping $\phi : \mathcal{V} \rightarrow [0, \infty]$ we have:

$$\mathbb{E}_{\omega \sim \Omega_n} (\phi(\psi_n(\omega))) \leq e n C_n \mathbb{E}_{A_0 \sim \mathcal{N}_{\mathbb{C}^{n \times n}}} \left(\frac{1}{n} \sum_{\lambda_0, v_0 : A_0 v_0 = \lambda_0 v_0} \phi(A_0, \lambda_0, v_0) \right). \quad (24)$$

Note that (24) can be understood as follows: let m_1 be the push-forward measure of ψ_n in \mathcal{V} and let m_2 be the measure in \mathcal{V} given by

$$m_2(X) = \mathbb{E}_{A_0 \sim \mathcal{N}_{\mathbb{C}^{n \times n}}} \left(\frac{1}{n} \# \{(\lambda_0, v_0) : A_0 v_0 = \lambda_0 v_0, (A_0, \lambda_0, v_0) \in X\} \right),$$

for any measurable set $X \subseteq \mathcal{V}$. Then, the Radon-Nikodim derivative dm_1/dm_2 is bounded above by enC_n .

Problem 2.34 Describe an alternative collection (Ω_n, ψ_n) which satisfies a sharper version of (24), with a constant in the place of nC_n . This would improve the running time of Algorithm `Random_initial_triple` below by a factor of $\mathcal{O}(nC_n)$.

We are now prepared to describe our random homotopy algorithm.

Algorithm 5 `Random_initial_triple`

Input: $A \in \mathbb{C}^{n \times n}$

Randomly choose $\omega \in \Omega_n$

$(A_0, \lambda_0, v_0) := \psi_n(\omega)$ (note that $v_0 = e_1$)

$(\zeta, w) := \text{Path-follow}(A, A_0, \lambda_0, v_0)$

Output: $(\zeta, w) \in (\mathbb{C} \times \mathbb{C}^n)^n$

Postconditions: The execution halts if the lifting of $\mathcal{L}_{A_0, A}$ at (λ_0, v_0) doesn't meet Σ' . In this case, (ζ, w) is an approximate eigenpair of A .

From Theorem 2.33, the expected running time of the computation of (A_0, λ_0, v_0) is $\mathcal{O}(n^3 C_n)$. Moreover, the expected number of homotopy steps in the execution of `Path-follow`(A, A_0, λ_0, v_0) is

$$S = \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}, \omega \sim \Omega_n} (K(A, \psi_n(\omega))) = \mathbb{E}_{\omega \sim \Omega_n} \left(\mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}} (K(A, \psi_n(\omega))) \right),$$

where $K(A, \psi_n(\omega))$ is as in Theorem 2.19. From (24), we have

$$\begin{aligned} S &\leq e n C_n \mathbb{E}_{A_0 \sim \mathcal{N}_{\mathbb{C}^{n \times n}}} \left(\frac{1}{n} \sum_{\lambda_0, v_0: A_0 v_0 = \lambda_0 v_0} \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}} (K(A, A_0, \lambda_0, v_0)) \right) \\ &\stackrel{\text{Th. 2.29}}{\leq} \mathcal{O}(n^3 C_n). \end{aligned}$$

We multiply the number of steps by $\mathcal{O}(n^3)$ to get the following complexity bound.

Theorem 2.35 *Algorithm `Random_initial_triple` returns (almost surely) an approximate eigenpair of its input $A \in \mathbb{C}^{n \times n}$. Its average cost satisfies*

$$\text{Avg_Cost}(n) = \mathcal{O}(n^6 C_n) \stackrel{\text{Lemma 10.1}}{\leq} \mathcal{O}(n^7). \quad \square$$

3 Some properties of the condition number μ

There is a general geometric framework for defining condition numbers, see [19, §14.3]. In our situation, this framework takes the following form.

If $(A, \lambda, v) \in \mathcal{W}$, then from the inverse function theorem the projection $\pi_1: \mathcal{V} \rightarrow \mathbb{C}^{n \times n}$ (cf. (3)), around (A, λ, v) , has a local inverse $\mathcal{U} \rightarrow \mathcal{V}, B \mapsto (B, G(B))$, that is defined on an open neighborhood \mathcal{U} of A in $\mathbb{C}^{n \times n}$. We call G the *solution map*. The map G decomposes as $G = (G_\lambda, G_v)$, where

$$G_\lambda : \mathcal{U} \rightarrow \mathbb{C} \quad \text{and} \quad G_v : \mathcal{U} \rightarrow \mathbb{P}(\mathbb{C}^n)$$

associate to matrices $B \in \mathcal{U}$ an eigenvalue and the corresponding eigenvector. Let

$$DG_\lambda(A) : \mathbb{C}^{n \times n} \rightarrow \mathbb{C} \quad \text{and} \quad DG_v(A) : \mathbb{C}^{n \times n} \rightarrow v^\perp$$

be the derivatives of these maps at A . The condition numbers for the eigenvalue λ and the eigenvector v of A are defined as follows:

$$\mu_\lambda(A, \lambda, v) := \|DG_\lambda(A)\| \quad \text{and} \quad \mu_v(A, \lambda, v) := \|DG_v(A)\|,$$

where the norms are the operator norms with respect to the chosen norms (on $\mathbb{C}^{n \times n}$ we use the Frobenius norm and on v^\perp the norm given by (1)).

Lemma 14.17 in [19] gives explicit descriptions of DG_λ and DG_v . Before stating it, we recall that if λ is an eigenvalue of A there exists $u \in \mathbb{P}(\mathbb{C}^n)$ (the *left eigenvector*) such that $(A - \lambda \text{Id})^* u = 0$. Recall the linear map $A_{\lambda, v} : v^\perp \rightarrow v^\perp$ introduced in (4).

Lemma 3.1 *Assume that $Av = \lambda v$ and λ has multiplicity 1. Then, the associated left eigenvector is*

$$u = v - i_{\mathbb{C}^n} A_{\lambda, v}^{-*} P_{v^\perp} A^* v. \quad (25)$$

Here we denoted $A_{\lambda, v}^{-*} := (A_{\lambda, v}^{-1})^*$. Note that $\langle u, v \rangle = \|v\|^2$.

PROOF. Take a representative such that $\|v\| = 1$ and let

$$z := i_{\mathbb{C}^n} A_{\lambda, v}^{-*} P_{v^\perp} A^* v \stackrel{(6)}{=} ((\text{Id} - vv^*)(A - \lambda \text{Id})^*(\text{Id} - vv^*))^\dagger A^* v.$$

From the definition of the Moore–Penrose pseudoinverse, z is the unique element in v^\perp that minimizes $\|(\text{Id} - vv^*)(A - \lambda \text{Id})^* z - A^* v\|$, that is we have $(\text{Id} - vv^*)(A - \lambda \text{Id})^* z = P_{v^\perp}(A^* v) = (\text{Id} - vv^*)A^* v$ or equivalently

$$(A - \lambda \text{Id})^* z = A^* v + tv \quad \text{for some } t \in \mathbb{C}.$$

Premultiplying both sides by v^* we have

$$v^*(A - \lambda \text{Id})^* z = v^* A^* v + t\|v\|^2 \Rightarrow 0 = (\lambda v)^* v + t\|v\|^2 = (\bar{\lambda} + t)\|v\|^2,$$

so we have $t = -\bar{\lambda}$ and then

$$(A - \lambda \text{Id})^*(v - z) = (A - \lambda \text{Id})^*v - A^*v + \bar{\lambda}v = 0,$$

that is, $v - z$ is a left eigenvector of A with associated (left) eigenvalue $\bar{\lambda}$ as wanted. \square

The following is Lemma 14.17 in [19].

Lemma 3.2 *Let $(A, \lambda, v) \in \mathcal{W}$ and let u be a left eigenvector of A with eigenvalue $\bar{\lambda}$. Then:*

- (a) *We have $\langle v, u \rangle \neq 0$.*
- (b) *The derivative of the solution map is given by $DG(A)(\dot{A}) = (\dot{\lambda}, \dot{v})$, where*

$$\dot{\lambda} = \frac{\langle \dot{A}v, u \rangle}{\langle v, u \rangle}, \quad \dot{v} = -A_{\lambda, v}^{-1} P_{v^\perp} \dot{A}v. \quad \square$$

The following result, which follows directly from Lemma 3.2, was already pointed out in [56] (see also [19, Prop. 14.15]).

Proposition 3.3 *Choosing the Frobenius norm on $T_A \mathbb{C}^{n \times n} = \mathbb{C}^{n \times n}$ and $\frac{1}{\|v\|} \|\cdot\|$ on v^\perp , the condition numbers μ_v for the eigenvector problem and μ_λ for the eigenvalue problem satisfy:*

$$\mu_\lambda(A, \lambda, v) = \|DG_\lambda(A)\| = \frac{\|u\| \|v\|}{|\langle u, v \rangle|} \stackrel{(25)}{=} \frac{\|u\|}{\|v\|} \leq \sqrt{1 + \mu^2(A, \lambda, v)}$$

and

$$\mu_v(A, \lambda, v) = \|DG_v(A)\| = \|A_{\lambda, v}^{-1}\| = \frac{\mu(A, \lambda, v)}{\|A\|_F}. \quad \square$$

PROOF OF PROPOSITION 2.7. The first two inequalities are immediate from Proposition 3.3. For the third one, note that

$$\|\dot{\Gamma}(t)\| = \|(\dot{A}, \dot{\lambda}, \dot{v})\| \leq \|\dot{A}\| \sqrt{1 + \mu_t^2 + (1 + \mu_t^2)} \leq \|\dot{A}\| \sqrt{6\mu_t^2}$$

the last inequality since $\mu_t \geq \frac{1}{\sqrt{2}}$ (Lemma 2.6). \square

Our last lemma is a version of Lemma 2.6 without the assumption that our point lies on \mathcal{W} .

Lemma 3.4 *For $A \in \mathbb{C}^{n \times n}$, $w \in \mathbb{C}^n$ and $\zeta \in \mathbb{C}$ with $|\zeta| \leq \|A\|_F$ we have*

$$\mu(A, \zeta, w) \geq \frac{1}{1 + \sqrt{1 - \frac{\|w^* A\|^2}{\|w\|^2 \|A\|_F^2}}} \geq \frac{1}{2}$$

PROOF. W.l.o.g we can assume that $w = e_1$ and write

$$A = \begin{pmatrix} \lambda & a^* \\ b & \hat{A} \end{pmatrix},$$

where $\lambda \in \mathbb{C}$ and $a, b \in \mathbb{C}^{n-1}$. Then, $A_{\zeta, w} \equiv \hat{A} - \zeta \mathbf{Id}_{n-1}$ and

$$\begin{aligned} \frac{\mu(A, \zeta, e_1)}{\|A\|_F} &= \|(\hat{A} - \zeta \mathbf{Id}_{n-1})^{-1}\| \geq \frac{1}{\|\hat{A} - \zeta \mathbf{Id}_{n-1}\|} \geq \frac{1}{\|\hat{A}\| + |\zeta|} \\ &\geq \frac{1}{\sqrt{\|A\|_F^2 - \|e_1^* A\|^2} + \|A\|_F}. \end{aligned}$$

The lemma follows. \square

4 Proofs of Theorem 2.8 and Corollary 2.9

It will be handy to use the definition of μ given in (10). We start with a very simple linear algebra lemma about the Moore-Penrose pseudoinverse.

Lemma 4.1 *Let $R, R' \in \mathbb{C}^{n \times n}$ be such that R has rank $n - 1$. Assume moreover that $\det(R') = 0$ and*

$$\|R - R'\| \leq \frac{\varepsilon}{\|R^\dagger\|}, \quad \text{for some } 0 \leq \varepsilon < 1.$$

Then, R' has rank $n - 1$ and

$$\frac{\|R^\dagger\|}{1 + \varepsilon} \leq \|R'^\dagger\| \leq \frac{\|R^\dagger\|}{1 - \varepsilon}.$$

PROOF. Let σ and σ' be the smallest nonzero singular values of R and R' , respectively. Note that $\sigma - \|R - R'\| \leq \sigma' \leq \sigma + \|R - R'\|$ (this is a classical fact proved for the first time in [58], see also [29, Cor. 8.6.2]). In particular, $\sigma' \geq \sigma - \|R - R'\| = \|R^\dagger\|^{-1} - \|R - R'\| > 0$, so R' has rank at least $n - 1$ and by hypothesis it has rank $n - 1$. Moreover, we have

$$\|R^\dagger\| = \frac{1}{\sigma} = \frac{1}{\sigma'} \frac{\sigma'}{\sigma} \leq \|R'^\dagger\| \frac{\sigma + \|R - R'\|}{\sigma} \leq (1 + \varepsilon) \|R'^\dagger\|.$$

The upper bound follows from a similar argument. \square

PROOF OF THEOREM 2.8. Choose representatives such that $\|v\| = \|v'\| = 1$ and let

$$Q := (\mathbf{Id} - vv^*)(A - \lambda \mathbf{Id})(\mathbf{Id} - vv^*), \quad Q' := (\mathbf{Id} - v'v'^*)(A' - \lambda' \mathbf{Id})(\mathbf{Id} - v'v'^*).$$

We have $\text{rank}(Q) = n - 1$ since $\mu(A, \lambda, v) < \infty$ by our assumption, cf. (10). We claim that

$$\|Q - Q'\| \leq \frac{\varepsilon}{\|Q^\dagger\|}, \quad (26)$$

which from Lemma 4.1 implies that Q' has rank $n - 1$ and

$$\frac{\|Q^\dagger\|}{1 + \varepsilon} \leq \|Q'^\dagger\| \leq \frac{\|Q^\dagger\|}{1 - \varepsilon},$$

that is (recall $\|A\|_F = \|A'\|_F = 1$),

$$\frac{\mu(A, \lambda, v)}{1 + \varepsilon} \leq \mu(A', \lambda', v') \leq \frac{\mu(A, \lambda, v)}{1 - \varepsilon},$$

as wanted.

It remains to prove the claim. To do so, let (A_t, λ_t, v_t) be a geodesic in $\mathbb{C}^{n \times n} \times \mathbb{C} \times \mathbb{P}(\mathbb{C}^n)$, parametrized by arc-length, joining (A, λ, v) and (A', λ', v') , so by hypothesis we have $t \in [0, \varepsilon/(4\sqrt{3}\|Q^\dagger\|)]$ and we can choose representatives v_t in such a way that

$$\dot{v}_t \perp v_t, \quad \|v_t\| = 1, \quad \|\dot{A}_t\|_F^2 + |\dot{\lambda}_t|^2 + \|\dot{v}_t\|^2 = 1, \quad \text{for all } t \in [0, \varepsilon/(4\sqrt{3}\|Q^\dagger\|)].$$

Let $Q_t := (\text{Id} - v_t v_t^*)(A_t - \lambda_t \text{Id})(\text{Id} - v_t v_t^*)$. In order to bound $\|\dot{Q}_t\|$ we first note that for $x \in \mathbb{C}^n$,

$$\begin{aligned} \|(\dot{v}_t v_t^* + v_t \dot{v}_t^*)(x)\| &= \|\dot{v}_t \langle x, v_t \rangle + v_t \langle x, \dot{v}_t \rangle\| = \sqrt{\|\dot{v}_t\|^2 |\langle x, v_t \rangle|^2 + \|v_t\|^2 |\langle x, \dot{v}_t \rangle|^2} \\ &= \|\dot{v}_t\| \sqrt{|\langle x, v_t \rangle|^2 + \left| \langle x, \frac{\dot{v}_t}{\|\dot{v}_t\|} \rangle \right|^2} \leq \|\dot{v}_t\| \|x\|, \end{aligned}$$

that is $\|\dot{v}_t v_t^* + v_t \dot{v}_t^*\| \leq \|\dot{v}_t\|$. On the other hand,

$$\|A_t - \lambda_t \text{Id}\| \leq \|A_t\| + |\lambda_t| \leq \|A_t\| + \|A_t\| \leq 2\|A_t\|_F = 2,$$

so computing the derivative of Q_t we see that

$$\begin{aligned} \|\dot{Q}_t\| &\leq 2\|\dot{v}_t v_t^* + v_t \dot{v}_t^*\| \|A_t - \lambda_t \text{Id}\| + \|\dot{A}_t - \dot{\lambda}_t \text{Id}\| \\ &\leq 4\|\dot{v}_t\| + \|\dot{A}_t - \dot{\lambda}_t \text{Id}\| \leq 4(\|\dot{v}_t\| + \|\dot{A}_t\| + |\dot{\lambda}_t|) \\ &\leq 4\sqrt{3}(\|\dot{v}_t\|^2 + \|\dot{A}_t\|^2 + |\dot{\lambda}_t|^2)^{1/2} = 4\sqrt{3}. \end{aligned}$$

Thus, we have $\|Q - Q'\| \leq \int_0^{\varepsilon/(4\sqrt{3}\|Q^\dagger\|)} 4\sqrt{3} dt \leq 4\sqrt{3}\varepsilon/(4\sqrt{3}\|Q^\dagger\|)$, finishing the proof of (26) and that of Theorem 2.8. \square

PROOF OF COROLLARY 2.9. Consider the portion $\mathcal{L}_{A, A'} = \{A_t\}$ of the great circle defined for $t \in [0, \alpha]$ where $\alpha = d_{\mathbb{S}}(A, A')$. Let (λ, v) be any eigenpair of A .

From the inverse function theorem, the map $t \mapsto \Gamma(t) = (A_t, \lambda_t, v_t) \in \mathcal{W}$ (given by the local inverse of π_1) is defined and, denoting $\text{dist}(t) := \text{dist}((A, \lambda, v), (A_t, \lambda_t, v_t))$, satisfies

$$\text{dist}(t) < \frac{\varepsilon}{4\sqrt{3}\mu(A, \lambda, v)}, \quad t \in [0, \delta) \quad (27)$$

for some maximal $\delta \in (0, \alpha]$. From Theorem 2.8 this implies

$$\mu(\Gamma(t)) \leq \frac{\mu(A, \lambda, v)}{1 - \varepsilon}, \quad t \in [0, \delta), \quad (28)$$

giving a global bound for the norm of the derivative of the solution map. Thus, $\Gamma(t)$ can be continuously extended to $[0, \delta]$ and from maximality of δ , either $\delta = \alpha$ or $\delta < \alpha$ and (27) must become an equality when changing t to δ . In this last case, which we will rule out by contradiction, note that from Proposition 2.7 and (28)

$$\text{dist}(\delta) \leq \int_0^\delta \|\dot{\Gamma}(t)\| dt \leq \sqrt{6} \int_0^\delta \mu(\Gamma(t)) dt \leq \sqrt{6} \delta \frac{\mu(A, \lambda, v)}{1 - \varepsilon}.$$

Now, in our range of values we have $\alpha = d_{\mathbb{S}}(A, A') = 2 \arcsin(\|A - A'\|_F/2) < \frac{1001}{1000} \|A - A'\|_F$ which, together with $\delta < \alpha$, implies

$$\text{dist}(\delta) < \frac{1001 \|A - A'\|_F \sqrt{6} \mu(A, \lambda, v)}{1000(1 - \varepsilon)} \leq \frac{1001 \varepsilon \sqrt{6}}{50000(1 - \varepsilon) \mu(A, \lambda, v)} < \frac{\varepsilon}{4\sqrt{3}\mu(A, \lambda, v)},$$

which is a contradiction (note that we have used the bound $\|A - A'\|_F \mu^2(A, \lambda, v) \leq \varepsilon/50$ and the fact that $\varepsilon < 1/2$). We thus conclude that $\delta = \alpha$ and the corollary follows from (27) and Theorem 2.8. \square

5 Condition-length homotopy continuation

The goal of this section is to fully describe the routine `Choose_step` (with which algorithm `Path-follow` will be complete) and to prove Theorem 2.19.

To do so, it will be useful to denote by $\beta(A, \zeta, w)$ the length (in the tangent space) of the Newton step with matrix $A \in \mathbb{S}$ and input $(\zeta, w) \in \mathbb{C} \times \mathbb{P}(\mathbb{C}^n)$. That is, if we take a representative such that $\|w\| = 1$,

$$\beta(A, \zeta, w)^2 := \left\| (DF_A(\zeta, w)|_{\mathbb{C} \times w^\perp})^{-1} F_A(\zeta, w) \right\|^2 = \left\| \begin{pmatrix} \dot{\lambda} \\ \dot{v} \end{pmatrix} \right\|^2 = |\dot{\lambda}|^2 + \|\dot{v}\|^2,$$

where $\dot{\lambda}, \dot{v}$ are given in (11). When $\|\dot{v}\|$ is small β approximates the length of the

Newton step also under the distance dist . Indeed,

$$\begin{aligned}\text{dist}_A((\zeta, w), (N_A(\zeta, w)))^2 &= |\dot{\lambda}|^2 + d_{\mathbb{P}}(w, w - \dot{v})^2 \\ &\stackrel{(2)}{=} |\dot{\lambda}|^2 + \left(\arccos \frac{|\langle w, w - \dot{v} \rangle|}{\|w\| \|w - \dot{v}\|} \right)^2 \\ &= |\dot{\lambda}|^2 + \left(\arccos \frac{1}{\sqrt{1 + \|\dot{v}\|^2}} \right)^2.\end{aligned}$$

It is easy to check that

$$\frac{9}{10} x^2 \leq \left(\arccos \frac{1}{\sqrt{1 + x^2}} \right)^2 \leq x^2 \quad \text{for } x \leq \frac{1}{3},$$

so we have proved that whenever $\|\dot{v}\| \leq 1/3$,

$$\frac{9}{10} \beta(A, \zeta, w) \leq \text{dist}_A((\zeta, w), (N_A(\zeta, w))) \leq \beta(A, \zeta, w). \quad (29)$$

(The upper bound inequality is valid regardless of the value of $\|\dot{v}\|$). The knowledge of β allows us to bound the distance from an approximate eigenpair to the associated exact eigenpair.

Lemma 5.1 *Assume that (ζ, w) is an approximate eigenpair of $A \in \mathbb{S}$ with associated eigenpair (λ, v) . Then, $\text{dist}_A((\zeta, w), (\lambda, v)) \leq 2\beta(A, \zeta, w)$. Moreover, if $\beta(A, \zeta, w) \leq 1/3$ we also have*

$$\frac{1}{2} \beta(A, \zeta, w) \leq \text{dist}_A((\zeta, w), (\lambda, v)) \leq 2\beta(A, \zeta, w).$$

PROOF. We have

$$\begin{aligned}\text{dist}_A((\zeta, w), (\lambda, v)) &\leq \text{dist}_A((\zeta, w), (N_A(\zeta, w))) + \text{dist}_A((N_A(\zeta, w)), (\lambda, v)) \\ &\stackrel{(29)}{\leq} \beta(A, \zeta, w) + \frac{1}{2} \text{dist}_A((\zeta, w), (\lambda, v)),\end{aligned}$$

the last from the definition of approximate eigenpair. The upper bound in the statement follows. Using a similar argument,

$$\begin{aligned}\text{dist}_A((\zeta, w), (\lambda, v)) &\geq \text{dist}_A((\zeta, w), (N_A(\zeta, w))) - \text{dist}_A((N_A(\zeta, w)), (\lambda, v)) \\ &\stackrel{(29)}{\geq} \frac{9}{10} \beta(A, \zeta, w) - \frac{1}{2} \text{dist}_A((\zeta, w), (\lambda, v)),\end{aligned}$$

and the lower bound follows as well. \square

5.1 The Lipschitz property of β

To describe `Choose_step` we need a set of constants satisfying a few relations. Not all of them are used in the description of `Choose_step`. Some only occur in the proof of the correctness of `Path-follow`. We will consider constants $c_1, c'_1, c_u, c'_u, c_*, c_4, c_5, c_6, c_7, K$. These are any collection of positive numbers satisfying the following:

$$\begin{aligned} \sqrt{3}c'_1 &\leq c_1 < \frac{1}{3}, & \sqrt{3}c'_u &\leq c_u - \frac{\frac{3}{2}c_1^2(\sqrt{3}-1)}{1-3c_1} \\ 4\sqrt{3}c_* &< 1, & c_4 &= c_* + (1+4\sqrt{3}c_*)(c_1+2c_u), & 4\sqrt{3}c_4 &< 1, \\ \frac{2(1+4\sqrt{3}c_*)}{1-4\sqrt{3}c_4}c_u &< Kc_* < \frac{1}{5}, & c_5 &= c'_u(1-4\sqrt{3}c_*) - 2\frac{2c_* + \frac{3}{2}c_1^2(1+4\sqrt{3}c_*)}{(1-3c_1)}, \\ c_6 &= \frac{c_5(1-3c_1) - 2(1+3c_1)c_*}{2(1+3c_1)(1+4\sqrt{3}c_4)}, & c_7 &= \min\left(\frac{c'_1(1-4\sqrt{3}c_*)}{(1+4\sqrt{3}c_4)}, c_6\right). \end{aligned}$$

A collection of parameters satisfying these constraints is shown in Table 1.

c'_1	10^{-3}	c_4	$0.005306\dots$
c_1	$\sqrt{3}c'_1$	c_5	$0.00099\dots$
c'_u	10^{-3}	c_6	$0.00038\dots$
c_u	$\sqrt{3}c'_u + \frac{3c_1^2(\sqrt{3}-1)}{2(1-3c_1)}$	K	64
c_*	10^{-4}	c_7	$0.00038\dots$

Table 1: Our choice of constants for `Path-follow`

In addition to these constants, for the sake of clarity, we spell out a working hypothesis that we will repeatedly use.

Hypothesis 5.2 We have $(A, \lambda, v) \in \mathcal{V}$, $\|A\|_F = 1$, and (ζ, w) an approximate eigenpair of A satisfying $|\zeta| \leq 1$ and

$$\mu(A, \lambda, v) \mathbf{dist}_A((\lambda, v), (\zeta, w)) < c_*.$$

Also, $\theta: [0, \pi) \rightarrow \mathbb{S}$, $\theta(s) = A_s$ is some arc-length parametrized half-great circle with $A_0 = A$.

It is easy to see from Theorem 2.12 that under hypotheses 5.2 we have $\beta(A, \zeta, w) < 1/3$ and in particular we can apply Lemma 5.1.

The main goal of this section is to prove the following result. Recall that we have defined

$$F_A(\lambda, v) = (A - \lambda \mathbf{Id})v.$$

Lemma 5.3 Under Hypothesis 5.2, let $s \leq c_1/\mu(A_0, \zeta, w)$ where $0 \leq c_1 < 1/3$. Let

$$\Phi = \left\| (DF_{A_0}(\zeta, w)|_{\mathbb{C} \times w^\perp})^{-1} \dot{A}_0 w \right\|.$$

Then,

$$\beta^-(s) \leq \beta(A_s, \zeta, w) \leq \beta^+(s),$$

where

$$\beta^-(s) = \frac{s\Phi - \beta(A_0, \zeta, w) - \frac{3}{2}c_1^2/\mu(A_0, \zeta, w)}{1 + 3c_1}, \quad \beta^+(s) = \frac{s\Phi + \beta(A_0, \zeta, w) + \frac{3}{2}c_1^2/\mu(A_0, \zeta, w)}{1 - 3c_1}.$$

For the proof we need some stepping stones.

Lemma 5.4 Let $A \in \mathbb{S}$, $\zeta \in \mathbb{C}$, $|\zeta| \leq 1$, $w \in \mathbb{C}^n$, $\|w\| = 1$. Then,

$$\left\| (DF_A(\zeta, w)|_{\mathbb{C} \times w^\perp})^{-1} \right\| \leq 3\mu(A, \zeta, w).$$

PROOF. This result is similar to [3, Proposition 6.6], but our assumptions are weaker and our definition of condition number is slightly different (see the proof of Theorem 2.12). We can assume that $w = e_1$ for the proof and write

$$A = \begin{pmatrix} \lambda & a^* \\ b & \hat{A} \end{pmatrix},$$

where $\lambda \in \mathbb{C}$ and $a, b \in \mathbb{C}^{n-1}$. Then, for $\dot{\zeta} \in \mathbb{C}$ and $\dot{w} \in \mathbb{C}^{n-1}$,

$$\begin{aligned} DF_A(\zeta, e_1) \begin{pmatrix} \dot{\zeta} \\ \dot{w} \end{pmatrix} &= (A - \zeta \mathbf{Id}) \begin{pmatrix} 0 \\ \dot{w} \end{pmatrix} - \dot{\zeta} e_1 \\ &= \begin{pmatrix} \lambda - \zeta & a^* \\ b & \hat{A} - \zeta \mathbf{Id}_{n-1} \end{pmatrix} \begin{pmatrix} 0 \\ \dot{w} \end{pmatrix} - \dot{\zeta} e_1 \\ &= \begin{pmatrix} -1 & a^* \\ 0 & \hat{A} - \zeta \mathbf{Id}_{n-1} \end{pmatrix} \begin{pmatrix} \dot{\zeta} \\ \dot{w} \end{pmatrix}. \end{aligned}$$

We thus have

$$\begin{aligned} \left\| (DF_A(\zeta, w)|_{\mathbb{C} \times w^\perp})^{-1} \right\| &= \left\| \begin{pmatrix} -1 & a^* \\ 0 & \hat{A} - \zeta \mathbf{Id}_{n-1} \end{pmatrix}^{-1} \right\| \\ &= \left\| \begin{pmatrix} -1 & a^*(\hat{A} - \zeta \mathbf{Id}_{n-1})^{-1} \\ 0 & (\hat{A} - \zeta \mathbf{Id}_{n-1})^{-1} \end{pmatrix} \right\| \\ &\leq \left\| \begin{pmatrix} 0 & 0 \\ 0 & (\hat{A} - \zeta \mathbf{Id}_{n-1})^{-1} \end{pmatrix} \right\| + \left\| \begin{pmatrix} -1 & a^*(\hat{A} - \zeta \mathbf{Id}_{n-1})^{-1} \\ 0 & 0 \end{pmatrix} \right\| \\ &\leq \|A_{\zeta, w}^{-1}\| + \sqrt{1 + \|a\|^2 \|A_{\zeta, w}^{-1}\|^2} \\ &= \mu(A, \zeta, w) + \sqrt{1 + \|a\|^2 \mu(A, \zeta, w)^2}, \end{aligned}$$

the last two claims in this chain by definition of $A_{\zeta,w}$ and the fact that $\|A\|_F = 1$. Now, from Lemma 3.4 we know that $\mu(A, \zeta, w) \geq (1 + \sqrt{1 - \|a\|^2})^{-1}$ which implies

$$\begin{aligned} \frac{\mu(A, \zeta, w) + \sqrt{1 + \|a\|^2 \mu(A, \zeta, w)^2}}{\mu(A, \zeta, w)} &= 1 + \sqrt{\|a\|^2 + \frac{1}{\mu(A, \zeta, w)^2}} \\ &\leq 1 + \sqrt{\|a\|^2 + (1 + \sqrt{1 - \|a\|^2})^2} \\ &= 1 + \sqrt{2 + 2\sqrt{1 - \|a\|^2}} \leq 3, \end{aligned}$$

the last inequality due to $0 \leq \|a\| \leq \|A\|_F = 1$. We have proved that

$$\left\| (DF_A(\zeta, w)|_{\mathbb{C} \times w^\perp})^{-1} \right\| \leq \mu(A, \zeta, w) + \sqrt{1 + \|a\|^2 \mu(A, \zeta, w)^2} \leq 3\mu(A, \zeta, w),$$

and the bound claimed in the lemma follows. \square

Lemma 5.5 *Under Hypothesis 5.2, let $s > 0$ satisfy $3\mu(A_0, \zeta, w) s < 1$. Then,*

$$\left\| \left[(DF_{A_s}(\zeta, w)|_{\mathbb{C} \times w^\perp})^{-1} (DF_{A_0}(\zeta, w)|_{\mathbb{C} \times w^\perp}) \right] \right\| \leq \frac{1}{1 - 3\mu(A_0, \zeta, w) s}$$

and

$$\left\| \left[(DF_{A_0}(\zeta, w)|_{\mathbb{C} \times w^\perp})^{-1} (DF_{A_s}(\zeta, w)|_{\mathbb{C} \times w^\perp}) \right] \right\| \leq 1 + 3\mu(A_0, \zeta, w) s.$$

PROOF. Note that

$$DF_{A_0}(\zeta, w)|_{\mathbb{C} \times w^\perp}(\dot{\eta}, \dot{x}) = DF_{A_s}(\zeta, w)|_{\mathbb{C} \times w^\perp}(\dot{\eta}, \dot{x}) + (A_0 - A_s)\dot{x},$$

hence

$$\begin{aligned} &\left\| \text{Id} - (DF_{A_0}(\zeta, w)|_{\mathbb{C} \times w^\perp})^{-1} (DF_{A_s}(\zeta, w)|_{\mathbb{C} \times w^\perp}) \right\| \\ &\leq \left\| (DF_{A_0}(\zeta, w)|_{\mathbb{C} \times w^\perp})^{-1} \right\| \|A_s - A_0\| \stackrel{\text{Lemma 5.4}}{\leq} 3\mu(A_0, \zeta, w) s \end{aligned}$$

where we used that $\|A_s - A_0\| \leq \|A_s - A_0\|_F \leq s$. The second claim in the statement is now obvious and the first one follows from the Banach lemma, $\|(\text{Id} + \Delta)^{-1}\| \leq (1 - \|\Delta\|)^{-1}$, valid for $\|\Delta\| < 1$. \square

Lemma 5.6 *Under Hypothesis 5.2, let $s > 0$ satisfy $3\mu(A_0, \zeta, w) s < 1$ and let*

$$\Phi_{\text{aux}}(s) = \left\| (DF_{A_0}(\zeta, w)|_{\mathbb{C} \times w^\perp})^{-1} (A_0 - A_s)w \right\|.$$

Then,

$$\frac{\Phi_{\text{aux}}(s) - \beta(A_0, \zeta, w)}{1 + 3\mu(A_0, \zeta, w) s} \leq \beta(A_s, \zeta, w) \leq \frac{\Phi_{\text{aux}}(s) + \beta(A_0, \zeta, w)}{1 - 3\mu(A_0, \zeta, w) s}.$$

PROOF. From the definition,

$$\begin{aligned}
\beta(A_s, \zeta, w) &= \left\| (DF_{A_s}(\zeta, w)|_{\mathbb{C} \times w^\perp})^{-1} F_{A_s}(\zeta, w) \right\| \\
&= \left\| \left[(DF_{A_s}(\zeta, w)|_{\mathbb{C} \times w^\perp})^{-1} (DF_{A_0}(\zeta, w)|_{\mathbb{C} \times w^\perp}) \right] \right. \\
&\quad \left. (DF_{A_0}(\zeta, w)|_{\mathbb{C} \times w^\perp})^{-1} (F_{A_0}(\zeta, w) + (A_s - A_0)w) \right\|. \tag{30}
\end{aligned}$$

Thus,

$$\beta(A_s, \zeta, w) \leq \left\| (DF_{A_s}(\zeta, w)|_{\mathbb{C} \times w^\perp})^{-1} (DF_{A_0}(\zeta, w)|_{\mathbb{C} \times w^\perp}) \right\| (\Phi_{\text{aux}}(s) + \beta(A_0, \zeta, w)).$$

Similarly, we have

$$\beta(A_s, \zeta, w) \geq \frac{\Phi_{\text{aux}}(s) - \beta(A_0, \zeta, w)}{\left\| (DF_{A_0}(\zeta, w)|_{\mathbb{C} \times w^\perp})^{-1} (DF_{A_s}(\zeta, w)|_{\mathbb{C} \times w^\perp}) \right\|}.$$

The statement now follows from Lemma 5.5. \square

PROOF OF LEMMA 5.3. We claim that we can write

$$A_s = A_0 + s\dot{A}_0 + \frac{s^2}{2}B,$$

for some B with $\|B\|_F \leq 1$. Indeed, this is an elementary observation which follows from the fact that A_s is a great circle in the sphere (w.l.o.g. one can choose A_s to be the circle parametrized by $(\cos s, \sin s, 0 \dots, 0)$ to prove it). As a consequence we have

$$\begin{aligned}
|\Phi_{\text{aux}}(s) - s\Phi| &\leq \left\| (DF_{A_0}(\zeta, w)|_{\mathbb{C} \times w^\perp})^{-1} (A_0 - A_s + s\dot{A}_0)w \right\| \\
&\leq \frac{s^2}{2} \left\| (DF_{A_0}(\zeta, w)|_{\mathbb{C} \times w^\perp})^{-1} \right\| \stackrel{\text{Lemma 5.4}}{\leq} \frac{3}{2}s^2\mu(A_0, \zeta, w).
\end{aligned}$$

The upper and lower bounds for $\beta(A_s, \zeta, w)$ in Lemma 5.3 now follow from this last estimate and Lemma 5.6. \square

5.2 The step's length in the homotopy continuation

The following result is crucial for the understanding of the homotopy algorithm. Its proof follows a logic which is similar to that of the proof of Corollary 2.9.

Proposition 5.7 *Under Hypothesis 5.2, let $s' > 0$ be any number such that*

$$c'_1 \leq \mu(A, \zeta, w)s' \leq c_1$$

and let s'' be any number such that

$$c'_u \leq \mu(A, \zeta, w)\beta^+(s'') \leq c_u,$$

where β^+ is as in Lemma 5.3. Let $\bar{s} = \min(s', s'')$. Then,

I. The (continuous) branch of the solution map $\pi_1^{-1} \circ \theta: [0, \bar{s}] \rightarrow \mathcal{V}$ with $\pi_1^{-1} \circ \theta(0) = (A, \lambda, v)$ is well defined.

II. For every $s \in [0, \bar{s}]$, let $(A_s, \lambda_s, v_s) := \pi_1^{-1} \circ \theta(s)$. Then,

$$\frac{1}{1 + 4\sqrt{3}c_4} \mu(A, \lambda, v) < \mu(A_s, \lambda_s, v_s) < \frac{\mu(A, \lambda, v)}{1 - 4\sqrt{3}c_4} \quad (31)$$

and

$$\text{dist}_{A_s}((\lambda_s, v_s), (\zeta, w)) < \frac{Kc_*}{\mu(A_s, \lambda_s, v_s)}. \quad (32)$$

In particular, (ζ, w) is an approximate eigenpair of A_s with associated eigenpair (λ_s, v_s) for every $s \in [0, \bar{s}]$.

Finally, the condition length $L_{\mu, 0, \bar{s}}(\pi_1^{-1}(A_s))$ (see (18)) of the curve $\pi_1^{-1}(A_s)$ is at least c_7 .

Before we prove Proposition 5.7 we make a few comments on how the proof of the proposition and Theorem 2.19 proceed. The hypotheses $\beta^+(s'')$ give us a bound on the distance from (ζ, w) to (λ_s, v_s) and ultimately (λ, v) to (λ_s, v_s) . Together with the bound on s' this allows us to invoke Theorem 2.12 and Theorem 2.8 to prove conclusions I and II of the proposition. The tricky part will be to prove the last statement that $L_{\mu, 0, \bar{s}}(\pi_1^{-1}(A_s)) \geq c_7$. This last statement gives us the upper bound on the number of steps in Theorem 2.19 as the condition length divided by c_7 . Now to see that

$$\int_0^{\bar{s}} \mu(A_s, \lambda_s, v_s) \|(\dot{A}_s, \dot{\lambda}_s, \dot{v}_s)\| ds = L_{\mu, 0, \bar{s}}(\pi_1^{-1}(A_s)) \geq c_7$$

it will suffice to prove that

$$\int_0^{\bar{s}} \|(\dot{A}_s, \dot{\lambda}_s, \dot{v}_s)\| ds \geq \frac{\text{constant}}{\mu(A, \lambda, v)},$$

since μ is almost constant on the interval. For this it will suffice to prove that one of

$$\int_0^{\bar{s}} \|\dot{A}_s\| ds \quad \text{or} \quad \int_0^{\bar{s}} \|\dot{v}_s\| ds$$

is greater than $\text{constant}/\mu(A, \lambda, v)$. The first integral is \bar{s} . We will see that if \bar{s} is small then $\text{dist}_A(v_s, v)$ is greater than or equal to some constant over $\mu(A, \lambda, v)$, and so is the integral of $\|\dot{v}_s\|$ which is the length of a path between v and v_s .

PROOF OF PROPOSITION 5.7. We prove the first part of the proposition. From the inverse function theorem and the continuity of μ , there exists a maximal $s^* \leq \bar{s}$ such that I and II hold changing $[0, \bar{s}]$ to $[0, s^*]$. The global upper bound for $\mu(\pi_1^{-1}(A_s))$ shown in (31) is in turn an upper bound for the derivative of the solution

map, for $s \in [0, s^*]$. A standard limit argument in compact sets then implies that $\pi_1^{-1} \circ \theta$ can be extended in a continuous manner to $[0, s^*]$, and because (31) and (32) are open conditions we must have one of the two following scenarios:

- i) $s_* = \bar{s}$ and both (31) and (32) hold changing s to s_* , or
- ii) At least one of (31) and (32) does not hold changing s to s_* .

We now discard the second option. Note that from Hypotheses 5.2 and Theorem 2.8,

$$\frac{1}{1 + 4\sqrt{3}c_*} \mu(A, \lambda, v) < \mu(A, \zeta, w) < \frac{\mu(A, \lambda, v)}{1 - 4\sqrt{3}c_*}. \quad (33)$$

Then, we have for every $s \in [0, s^*]$

$$\text{dist}((A, \zeta, w), (A_s, \zeta, w)) = \|A - A_s\|_F < s^* \leq s' \leq \frac{c_1}{\mu(A, \zeta, w)} \leq \frac{(1 + 4\sqrt{3}c_*)c_1}{\mu(A, \lambda, v)},$$

and because $s^* \leq s''$, from Lemma 5.3 we have

$$\beta(A_s, \zeta, w) \leq \beta^+(s) \leq \frac{c_u}{\mu(A, \zeta, w)} < \frac{(1 + 4\sqrt{3}c_*)c_u}{\mu(A, \lambda, v)}.$$

From Lemma 5.1 (recall that from II and Theorem 2.12, (ζ, w) is an approximate eigenpair of A_s with associated eigenpair (λ_s, v_s) for $s \leq s_*$), this last inequality implies

$$\text{dist}_{A_s}((\zeta, w), (\lambda_s, v_s)) \leq 2\beta(A_s, \zeta, w) < \frac{2(1 + 4\sqrt{3}c_*)c_u}{\mu(A, \lambda, v)}. \quad (34)$$

We have thus proved that for every $s \in [0, s^*]$,

$$\begin{aligned} \text{dist}((A, \lambda, v), (A_s, \lambda_s, v_s)) &\leq \text{dist}_A((\lambda, v), (\zeta, w)) + \text{dist}((A, \zeta, w), (A_s, \zeta, w)) \\ &\quad + \text{dist}_{A_s}((\zeta, w), (\lambda_s, v_s)) \\ &< \frac{c_*}{\mu(A, \lambda, v)} + \frac{(1 + 4\sqrt{3}c_*)c_1}{\mu(A, \lambda, v)} + \frac{2(1 + 4\sqrt{3}c_*)c_u}{\mu(A, \lambda, v)} \\ &= \frac{c_4}{\mu(A, \lambda, v)}. \end{aligned}$$

Then, from Theorem 2.8 (note the strict inequality in the displayed formula above),

$$\frac{1}{1 + 4\sqrt{3}c_4} \mu(A, \lambda, v) < \mu(A_{s^*}, \lambda_{s^*}, v_{s^*}) < \frac{\mu(A, \lambda, v)}{1 - 4\sqrt{3}c_4}. \quad (35)$$

Thus, (31) holds at s^* and moreover

$$\text{dist}_{A_{s^*}}((\lambda_{s^*}, v_{s^*}), (\zeta, w)) \underset{(34),(35)}{<} \frac{2(1 + 4\sqrt{3}c_*)c_u}{(1 + 4\sqrt{3}c_4)\mu(A_{s^*}, \lambda_{s^*}, v_{s^*})} < \frac{Kc_*}{\mu(A_{s^*}, \lambda_{s^*}, v_{s^*})}.$$

That is, (32) holds at s^* and we can discard option ii), proving the first part of the proposition.

For the last claim of the proposition, note that the condition length of the curve $\pi_1^{-1}(A_s)$ is

$$\begin{aligned} L_{\mu,0,\bar{s}}(\pi_1^{-1}(A_s)) &= \int_0^{\bar{s}} \mu(\pi_1^{-1}(A_s)) \left\| D\pi_1^{-1}(A_s) \dot{A}_s \right\| ds \\ &\stackrel{(31)}{\geq} \frac{\mu(A, \lambda, v)}{1 + 4\sqrt{3}c_4} \int_0^{\bar{s}} \left\| D\pi_1^{-1}(A_s) \dot{A}_s \right\| ds. \end{aligned}$$

If $\bar{s} = s'$ then using that the integrand is greater than or equal to 1 and (33) we have

$$L_{\mu,0,\bar{s}}(\pi_1^{-1}(A_s)) \geq \frac{\mu(A, \lambda, v)}{1 + 4\sqrt{3}c_4} s' \geq \frac{c'_1}{1 + 4\sqrt{3}c_4} \frac{\mu(A, \lambda, v)}{\mu(A, \zeta, w)} \geq \frac{c'_1(1 - 4\sqrt{3}c_*)}{(1 + 4\sqrt{3}c_4)}.$$

Assume now that $\bar{s} = s''$. Then we have

$$\begin{aligned} L_{\mu,0,\bar{s}}(\pi_1^{-1}(A_s)) &\geq \frac{\mu(A, \lambda, v)}{1 + 4\sqrt{3}c_4} \int_0^{\bar{s}} \left\| D\pi_1^{-1}(A_s) \dot{A}_s \right\| ds \\ &\geq \frac{\mu(A, \lambda, v)}{1 + 4\sqrt{3}c_4} \text{dist}((A_0, \lambda_0, v_0), (A_{\bar{s}}, \lambda_{\bar{s}}, v_{\bar{s}})). \end{aligned}$$

Now, note that (recall $(A_0, \lambda_0, v_0) = (A, \lambda, v)$)

$$\begin{aligned} \text{dist}((A_0, \lambda_0, v_0), (A_{\bar{s}}, \lambda_{\bar{s}}, v_{\bar{s}})) &\geq \text{dist}((A_0, \zeta, w), (A_{\bar{s}}, \lambda_{\bar{s}}, v_{\bar{s}})) - \text{dist}_{A_0}((\lambda_0, v_0), (\zeta, w)) \\ &\geq \text{dist}_{A_{\bar{s}}}((\zeta, w), (\lambda_{\bar{s}}, v_{\bar{s}})) - \frac{c_*}{\mu(A, \lambda, v)}. \end{aligned}$$

We need a lower bound for this last term. We first note that from Lemma 5.1 and Hypothesis 5.2,

$$\beta(A, \zeta, w) \leq \frac{2c_*}{\mu(A, \lambda, v)} \leq \frac{2c_*}{(1 - 4\sqrt{3}c_*)\mu(A, \zeta, w)}, \quad (36)$$

the last by (33). Using this last bound and, again, Lemmas 5.1 and 5.3, we have

$$\begin{aligned} 2 \frac{1 + 3c_1}{1 - 3c_1} \text{dist}_{A_{\bar{s}}}((\zeta, w), (\lambda_{\bar{s}}, v_{\bar{s}})) &\geq \frac{1 + 3c_1}{1 - 3c_1} \beta(A_{\bar{s}}, \zeta, w) \\ &\geq \frac{1 + 3c_1}{1 - 3c_1} \beta^-(\bar{s}) = \beta^+(\bar{s}) - 2 \frac{\beta(A, \zeta, w) + \frac{3}{2}c_1^2/\mu(A, \zeta, w)}{1 - 3c_1} \\ &\geq \frac{c'_u}{\mu(A, \zeta, w)} - 2 \frac{\beta(A, \zeta, w) + \frac{3}{2}c_1^2/\mu(A, \zeta, w)}{1 - 3c_1} \\ &\geq \frac{c'_u(1 - 4\sqrt{3}c_*)}{\mu(A, \lambda, v)} - 2 \frac{2c_* + \frac{3}{2}c_1^2(1 + 4\sqrt{3}c_*)}{(1 - 3c_1)\mu(A, \lambda, v)} = \frac{c_5}{\mu(A, \lambda, v)}. \end{aligned}$$

For the inequality in the third line we used the assumption in the statement and $s = s''$. For the inequality in the fourth line, we used (33) and (36). For the equality in the fourth line we used the definition of c_5 .

We have thus shown that if $\bar{s} = s''$, then

$$\begin{aligned} L_{\mu,0,\bar{s}}(\pi_1^{-1}(A_s)) &\geq \frac{\mu(A,\lambda,v)}{1+4\sqrt{3}c_4} \left(\frac{c_5(1-3c_1)}{2(1+3c_1)\mu(A,\lambda,v)} - \frac{c_*}{\mu(A,\lambda,v)} \right) \\ &= \frac{c_5(1-3c_1) - 2(1+3c_1)c_*}{2(1+3c_1)(1+4\sqrt{3}c_4)} = c_6. \end{aligned}$$

Since $c_7 \leq c_6$ the proof is complete. \square

We can finally describe the subroutine `Choose_step`. It is important to note that given any matrix $A \in \mathbb{C}^{n \times n}$, one can compute in $\mathcal{O}(n^3)$ arithmetic operations, for example by first reducing A to tridiagonal Hessenberg form and then using the main result of [31], a number r such that $\|A\| \leq r \leq \sqrt{3}\|A\|$. That is, we can compute operator norms within a factor of $\sqrt{3}$ and, consequently, we can compute μ within a factor of $\sqrt{3}$.

Algorithm 6 `Choose_step`

Input: $B, \dot{A} \in \mathbb{S}$, and $(\zeta, w) \in \mathbb{C} \times \mathbb{C}^n$, $\|w\| = 1$

compute $r > 0$ such that $\mu(B, \zeta, w) \leq r \leq \sqrt{3}\mu(B, \zeta, w)$

$s' := c_1/r$

$\Phi := \left\| (DF_B(\zeta, w)|_{\mathbb{C} \times w^\perp})^{-1} \dot{A} w \right\|$

compute s'' , the solution of the following linear equation

$$\frac{\Phi s'' + \beta(B, \zeta, w) + \frac{3}{2}c_1^2\sqrt{3}/r}{1-3c_1} = \frac{c_u}{r}$$

$\bar{s} := \min(s', s'')$

Output: $\Delta s = \bar{s} \in [0, \pi]$

The step size computed by `Choose_step` cannot be too small, as we show now.

Proposition 5.8 *The value Δs returned by `Choose_step`(B, \dot{A}, ζ, w) satisfies*

$$\Delta s \geq \frac{R}{\mu^2(B, \zeta, w)}$$

with $R = c_7(1-4\sqrt{3}c_4)^2/6 > 0$.

5.3 Proof of Theorem 2.19 and Proposition 5.8

We now prove Theorem 2.19, and the proof of Proposition 5.8 will follow straightforward from our arguments.

From the definition of Path-follow it is clear that we can assume that $\|A_0\|_F = \|A\|_F = 1$. We further assume that the constants $c'_1, c_1, c'_u, c_u c_*, c_4, c_5, c_6, c_7$ and K take the values in Table 1 and denote by $\pi_1^{-1}(\mathcal{L}_{A_0, A})$ the lift of $\mathcal{L}_{A_0, A}$ with origin (A_0, λ_0, v_0) . Note that this lift is well-defined since, by hypothesis, $\mathcal{L}_{A_0, A} \cap \Sigma = \emptyset$.

Let B_i be the matrix B at the beginning of the i th iteration of Path-follow. Also, let (λ_i, v_i) be such that (B_i, λ_i, v_i) is the (unique) triple in $\pi_1^{-1}(\mathcal{L}_{A_0, A})$ above B_i .

We first prove that for all $i \geq 0$, (ζ_i, w_i) is an approximate zero of B_i with associated eigenpair (λ_i, v_i) and satisfies

$$\text{dist}_{B_i}((\zeta_i, w_i), (\lambda_i, v_i)) < \frac{c_*}{\mu(B_i, \lambda_i, v_i)}.$$

We reason by induction. The step $i = 0$ is true by hypothesis (recall Definition 2.17). For the induction step, note that the s' defined by Choose_step satisfies (we omit the subindices i in (B_i, ζ_i, w_i) in the next few lines)

$$\frac{c'_1}{\mu(B, \zeta, w)} \leq \frac{c_1}{\sqrt{3}\mu(B, \zeta, w)} \leq s' \leq \frac{c_1}{\mu(B, \zeta, w)}.$$

Moreover, s'' satisfies

$$\begin{aligned} \beta^+(s'') &= \frac{\Phi s'' + \beta(B, \zeta, w) + \frac{3}{2}c_1^2/\mu(B, \zeta, w)}{1 - 3c_1} \\ &\leq \frac{\Phi s'' + \beta(B, \zeta, w) + \frac{3}{2}c_1^2\sqrt{3}/r}{1 - 3c_1} = \frac{c_u}{r} \leq \frac{c_u}{\mu(B, \zeta, w)}, \end{aligned}$$

and

$$\begin{aligned} \beta^+(s'') &= \frac{\Phi s'' + \beta(B, \zeta, w) + \frac{3}{2}c_1^2/\mu(B, \zeta, w)}{1 - 3c_1} \geq \frac{\Phi s'' + \beta(B, \zeta, w) + \frac{3}{2}c_1^2/r}{1 - 3c_1} \\ &= \frac{c_u - \frac{3}{2}c_1^2(\sqrt{3} - 1)/(1 - 3c_1)}{r} \geq \frac{c_u - \frac{3}{2}c_1^2(\sqrt{3} - 1)/(1 - 3c_1)}{\sqrt{3}\mu(B, \zeta, w)} \geq \frac{c'_u}{\mu(B, \zeta, w)}. \end{aligned}$$

We are thus under the hypothesis of Proposition 5.7 with $\bar{s} = \Delta s$ (by construction in Choose_step), which guarantees that (ζ_i, w_i) is an approximate eigenpair of the matrix B_{i+1} . Moreover, Proposition 5.7 also implies

$$\text{dist}_{B_{i+1}}((\zeta_i, w_i), (\lambda_{i+1}, v_{i+1})) < \frac{Kc_*}{\mu(B_{i+1}, \lambda_{i+1}, v_{i+1})}.$$

Since $Kc_* < 1/5$ we deduce (using Theorem 2.12) that (ζ_i, w_i) is an approximate eigenpair of B_{i+1} with associated eigenpair (λ_{i+1}, v_{i+1}) . Consequently, after three

steps of Newton iteration, we have that (ζ_{i+1}, w_{i+1}) satisfies (recall, $K = 64$)

$$\begin{aligned} \mathsf{dist}_{B_{i+1}}((\zeta_{i+1}, w_{i+1}), (\lambda_{i+1}, v_{i+1})) &< \frac{1}{2^{2^3-1}} \frac{64c_*}{\mu(B_{i+1}, \lambda_{i+1}, v_{i+1})} \\ &= \frac{c_*}{2\mu(B_{i+1}, \lambda_{i+1}, v_{i+1})}. \end{aligned} \quad (37)$$

If $|\zeta_{i+1}| \leq 1$, this finishes the induction step. Otherwise, the algorithm divides ζ_{i+1} by its norm, in that case we have

$$\begin{aligned} \mathsf{dist}_{B_{i+1}}\left(\left(\frac{\zeta_{i+1}}{|\zeta_{i+1}|}, w_{i+1}\right), (\lambda_{i+1}, v_{i+1})\right) &\leq \mathsf{dist}_{B_{i+1}}\left(\left(\frac{\zeta_{i+1}}{|\zeta_{i+1}|}, w_{i+1}\right), (\zeta_{i+1}, w_{i+1})\right) + \\ &\quad \mathsf{dist}_{B_{i+1}}((\zeta_{i+1}, w_{i+1}), (\lambda_{i+1}, v_{i+1})) \\ &\leq |\zeta_{i+1}| - 1 + \frac{c_*}{2\mu(B_{i+1}, \lambda_{i+1}, v_{i+1})} \end{aligned}$$

On the other hand, from (37) we have (use $|\lambda_{i+1}| \leq \|B_{i+1}\|_F = 1$)

$$|\zeta_{i+1} - \lambda_{i+1}| < \frac{c_*}{2\mu(B_{i+1}, \lambda_{i+1}, v_{i+1})} \Rightarrow |\zeta_{i+1}| < 1 + \frac{c_*}{2\mu(B_{i+1}, \lambda_{i+1}, v_{i+1})},$$

so we have

$$\mathsf{dist}_{B_{i+1}}\left(\left(\frac{\zeta_{i+1}}{|\zeta_{i+1}|}, w_{i+1}\right), (\lambda_{i+1}, v_{i+1})\right) < \frac{c_*}{\mu(B_{i+1}, \lambda_{i+1}, v_{i+1})},$$

and the induction step is finished also in the case that $|\zeta_{i+1}| > 1$.

The induction step is complete. In particular, this shows the last part of the statement.

To show the complexity bounds, assume **Path-follow** has performed $q+\ell$ iterations and let $0 = s_0 < s_1 < \dots < s_q < \dots < s_{q+\ell}$ be the corresponding values of s . Then we have

$$\begin{aligned} L_{\mu, s_q, s_{q+\ell}}(\pi_1^{-1}(B_s)) &= \sum_{i=1}^{\ell} \int_{s_{q+i-1}}^{s_{q+i}} \mu(A_s, \lambda_s, v_s) \|(\dot{A}_s, \dot{\lambda}_s, \dot{v}_s)\| ds \\ &= \sum_{i=1}^{\ell} L_{\mu, s_{q+i-1}, s_{q+i}}(\pi_1^{-1}(A_s)) \geq \ell c_7 \end{aligned}$$

the inequality by the last claim of Proposition 5.7. But the algorithm halts as soon as $s_{q+\ell} = \alpha$, i.e., as soon as

$$L_{\mu, s_q, s_{q+\ell}}(\pi_1^{-1}(A_s)) = L_{\mu, s_q, \alpha}(\pi_1^{-1}(A_s)),$$

which occurs as soon as $\ell \geq c_7^{-1} \int_{s_q}^{\alpha} \mu(A_s, \lambda_s, v_s) \|(\dot{A}_s, \dot{\lambda}_s, \dot{v}_s)\| ds$, as claimed in the theorem (note that $C := c_7^{-1} \leq 3000$).

We finally prove Proposition 5.8. From Proposition 5.7 we have

$$\begin{aligned} \mu^2(B, \zeta, w) \Delta s &\stackrel{(31)}{\geq} (1 - 4\sqrt{3}c_4)^2 \int_0^{\Delta s} \mu^2(A_s, \lambda_s, v_s) ds \\ &\stackrel{\text{Prop. 2.7}}{\geq} \frac{(1 - 4\sqrt{3}c_4)^2}{6} L_{\mu, 0, \Delta s}(\pi_1^{-1}(A_s)) \geq \frac{c_7(1 - 4\sqrt{3}c_4)^2}{6}, \end{aligned}$$

so Proposition 5.8 follows. \square

6 Integration in the solution variety

6.1 The coarea formula

On a Riemannian manifold M there is a well-defined measure vol_M obtained by integrating the indicator functions $\mathbb{1}_A$ of Borel-measurable subsets $A \subseteq M$ against the volume form dM of M ,

$$\text{vol}_M(A) := \int_M \mathbb{1}_A dM.$$

Dividing $\mathbb{1}_A$ by $\text{vol}_M(M)$ if $\text{vol}_M(M) < \infty$, this leads to a natural notion of *uniform distribution* on M , which we will denote by $\mathcal{U}(M)$. More generally, we will call any measurable function $f: M \rightarrow [0, \infty]$ such that $\int_M f dM = 1$ a *probability density* on M . We abuse notation and sometimes denote $\mathbb{E}_{x \in M}$ instead of $\mathbb{E}_{x \sim \mathcal{U}(M)}$.

The coarea formula (a modern classical formula due to Federer [28], see the Appendix of [30] for a smooth version) is an extension of the transformation formula to not necessarily bijective smooth maps between Riemannian manifolds. In order to state it, we first need to generalize the notion of Jacobians.

Suppose that M, N are Riemannian manifolds of dimensions m, n , respectively such that $m \geq n$. Let $\psi: M \rightarrow N$ be a smooth map. By definition, the derivative $D\psi(x): T_x M \rightarrow T_{\psi(x)} N$ at a regular point $x \in M$ is surjective. Hence the restriction of $D\psi(x)$ to the orthogonal complement of its kernel yields a linear isomorphism. The absolute value of its determinant is called the *normal Jacobian* (sometimes called *normal determinant* in the context of Linear Algebra, see [1]) of ψ at x and is denoted by $\text{NJ}\psi(x)$. We set $\text{NJ}\psi(x) := 0$ if x is not a regular point.

If y is a regular value of ψ , then the fiber $F_y := \psi^{-1}(y)$ is a Riemannian submanifold of M of dimension $m - n$, and it makes sense to integrate functions on F_y . Moreover, Sard's lemma states that almost all $y \in N$ are regular values.

We can now state the *coarea formula*.

Theorem 6.1 (Coarea formula) *Suppose that M, N are Riemannian manifolds of dimensions m, n , respectively, and let $\psi: M \rightarrow N$ be a surjective smooth map*

such that $D\psi$ is surjective a.e. Put $F_y = \psi^{-1}(y)$. Then we have for any function $\chi: M \rightarrow \mathbb{R}$ that is integrable with respect to the volume measure of M that

$$\int_M \chi \, dM = \int_{y \in N} \left(\int_{F_y} \frac{\chi}{\text{NJ}\psi} \, dF_y \right) dN,$$

and the integrals involved are well-defined. \square

It should be clear that this result contains the change of variables formula as a special case. Moreover, if we apply the coarea formula to the projection $\pi_2: M \times N \rightarrow N$, $(x, y) \mapsto y$, we retrieve Fubini's theorem since $\text{NJ}\pi_2 = 1$.

6.2 Coarea formula and double fibrations

The coarea formula can be readily applied to the following situation. Assume that three Riemannian manifolds M , N_1 , N_2 are equipped with surjective smooth mappings $\pi_1: M \rightarrow N_1$ and $\pi_2: M \rightarrow N_2$ whose derivatives are a.e. surjective, so $\text{NJ}\pi_1$ and $\text{NJ}\pi_2$ are a.e. nonzero. Let $\chi: M \rightarrow [0, \infty)$ be a measurable mapping. From Theorem 6.1 applied to π_1 we have (here dx and dy stand for the volume forms in M and N_1 , respectively)

$$\int_{x \in M} \chi(x) \text{NJ}\pi_1(x) \, dx = \int_{y \in N_1} \int_{x \in \pi_1^{-1}(y)} \chi(x) \, dx \, dy.$$

On the other hand, Theorem 6.1 applied to π_2 yields

$$\int_{x \in M} \chi(x) \text{NJ}\pi_1(x) \, dx = \int_{z \in N_2} \int_{x \in \pi_2^{-1}(z)} \frac{\text{NJ}\pi_1(x)}{\text{NJ}\pi_2(x)} \chi(x) \, dx \, dz.$$

We thus have the following result.

Theorem 6.2 *Let M , N_1 , N_2 be Riemannian manifolds equipped with surjective smooth mappings $\pi_1: M \rightarrow N_1$ and $\pi_2: M \rightarrow N_2$ whose derivatives are a.e. surjective. Let $\chi: M \rightarrow [0, \infty)$ be a measurable mapping. Then,*

$$\int_{y \in N_1} \int_{x \in \pi_1^{-1}(y)} \chi(x) \, dx \, dy = \int_{z \in N_2} \int_{x \in \pi_2^{-1}(z)} \frac{\text{NJ}\pi_1(x)}{\text{NJ}\pi_2(x)} \chi(x) \, dx \, dz.$$

(Note that when M and N_1 have the same dimension, one can replace the inner integral in the left-hand side by a (finite or enumerable) sum). \square

In the next sections we shall apply this result in two different contexts. A linear algebra argument simplifies the computation of the quotient of Normal Jacobians. Let E and F be finite dimensional, complex Euclidean vector spaces and let $\varphi: E \rightarrow$

F be a surjective linear mapping. Consider the graph $\Gamma := \{(x, \varphi(x)) \mid x \in E\}$ of φ . Then, Γ is a linear subspace of $E \times F$ and the two projections

$$p_1: \Gamma \rightarrow E, (x, \varphi(x)) \mapsto x, \quad p_2: \Gamma \rightarrow F, (x, \varphi(x)) \mapsto \varphi(x)$$

are linear maps. Note that p_1 is an isomorphism and p_2 is surjective as φ is.

Lemma 6.3 *Under the above assumptions, we have*

$$\frac{\text{NJ}p_1}{\text{NJ}p_2} = |\det(\varphi\varphi^*)|^{-1}.$$

PROOF. This result is [14, Lemma 3b), p. 242], although we rewrite it for complex vector spaces here (note the comment in the proof of [14, Theorem 5, p. 243]). \square

6.3 The solution variety for the eigenpair problem

Recall from §2.2, we have the two projections

$$\pi_1: \mathcal{V} \rightarrow \mathbb{C}^{n \times n}, (A, \lambda, v) \mapsto A \quad \text{and} \quad \pi_2: \mathcal{V} \rightarrow \mathbb{C} \times \mathbb{P}(\mathbb{C}^n), (A, \lambda, v) \mapsto (\lambda, v)$$

and, for $(A, \lambda, v) \in \mathcal{V}$, the linear operator $A_{\lambda, v}: v^\perp \rightarrow v^\perp$ given by $P_{v^\perp}(A - \lambda \mathbf{Id})|_{v^\perp}$. In order to apply Theorem 6.2 we first need to compute the quotient of normal Jacobians there.

Proposition 6.4 *Let $p := (A, \lambda, v) \in \mathcal{W}$ and choose a representative such that $\|v\| = 1$. Then, the derivative $D\pi_1(p): T_p \mathcal{W} \rightarrow T_A \mathbb{C}^{n \times n}$ is an isomorphism, the derivative $D\pi_2(p): T_p \mathcal{W} \rightarrow T_{(\lambda, v)}(\mathbb{C} \times \mathbb{P}(\mathbb{C}^n))$ is surjective, and we have*

$$\frac{\text{NJ}\pi_1(p)}{\text{NJ}\pi_2(p)} = |\det A_{\lambda, v}|^2 = \det(A_{\lambda, v} A_{\lambda, v}^*).$$

PROOF. By unitary invariance, we may assume without loss of generality that $v = e_1 = (1, 0, \dots, 0)$, and then

$$A = \begin{pmatrix} \lambda & c^* \\ 0 & B \end{pmatrix}, \quad c \in \mathbb{C}^{n-1}, B \in \mathbb{C}^{(n-1) \times (n-1)},$$

so $A_{\lambda, v} = B - \lambda \mathbf{Id}_{n-1}$. Let $\Gamma = \{(\dot{A}, DG(A)\dot{A}) : \dot{A} \in \mathbb{C}^{n \times n}\} \subseteq \mathbb{C}^{n \times n} \times \mathbb{C} \times v^\perp$ where G is the appropriate branch of the solution map defined in some open neighborhood of A . We are under the hypotheses of Lemma 6.3 so we have

$$\frac{\text{NJ}\pi_1(p)}{\text{NJ}\pi_2(p)} = \frac{\text{NJ}(D\pi_1(p))}{\text{NJ}(D\pi_2(p))} = \det(DG(A, \lambda, v)DG(A, \lambda, v)^*)^{-1}.$$

From Lemmas 3.1 and 3.2, we have

$$DG(A, \lambda, v)\dot{A} = \begin{pmatrix} \langle \dot{A}v, v - i_{\mathbb{C}^n} A_{\lambda, v}^{-*} P_{v^\perp} A^* v \rangle \\ -A_{\lambda, v}^{-1} P_{v^\perp} \dot{A}v \end{pmatrix} = \begin{pmatrix} v^* - v^* A i_{\mathbb{C}^n} A_{\lambda, v}^{-1} P_{v^\perp} \\ -A_{\lambda, v}^{-1} P_{v^\perp} \end{pmatrix} \dot{A}v = R \dot{A}v,$$

where, in the last displayed formula, we are denoting by $R: \mathbb{C}^n \rightarrow \mathbb{C} \times v^\perp \equiv \mathbb{C}^n$ the linear operator multiplying by $\dot{A}v$ in the previous formula. A standard linear algebra argument then shows that $\det(DG(A, \lambda, v)DG(A, \lambda, v)^*) = \det(RR^*) = |\det(R)|^2$. Now, we can identify

$$i_{\mathbb{C}^n} \equiv \begin{pmatrix} 0 \\ \text{Id}_{n-1} \end{pmatrix}, \quad P_{v^\perp} \equiv (0 \quad \text{Id}_{n-1}),$$

which implies that in standard basis we have

$$R = \begin{pmatrix} 1 & * \\ 0 & -A_{\lambda, v}^{-1} \end{pmatrix},$$

thus showing that $|\det(R)|^2 = |\det A_{\lambda, v}^{-1}|^2$, and the proposition follows. \square

We are now ready to rewrite Theorem 6.2 in this setting. The following is an important technical result that we will use several times.

Proposition 6.5 *Let $\chi: \mathcal{V} \rightarrow [0, \infty)$ be a measurable mapping. Then,*

$$\int_{A \in \mathbb{C}^{n \times n}} \sum_{\lambda, v: Av = \lambda v} \chi(A, \lambda, v) dA = \int_{(\lambda, v) \in \mathbb{C} \times \mathbb{P}(\mathbb{C}^n)} \int_{A: Av = \lambda v} \chi(A, \lambda, v) |\det(A_{\lambda, v})|^2 dA d(\lambda, v).$$

Moreover, assume that χ is unitarily invariant in the sense that $\chi(A, \lambda, v) = \chi(UAU^*, \lambda, Uv)$ for any unitary matrix $U \in \mathcal{U}_n$. Fix any a.e. continuous mapping $\mathbb{C}^n \setminus \{0\} \rightarrow \mathcal{U}_n$, $v \mapsto U_v$ such that $U_v e_1 = v/\|v\|$ for all v . Then, for every $\hat{A} \in \mathbb{C}^{n \times n}$ and $\sigma > 0$,

$$\begin{aligned} & \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}(\hat{A}, \sigma^2)} \left(\sum_{\lambda, v: Av = \lambda v} \chi(A, \lambda, v) \right) \\ &= \frac{1}{\Gamma(n)\sigma^{2(n-1)}} \mathbb{E}_v \mathbb{E}_{(\lambda, w, B)} \left(e^{-\frac{\|\hat{y}_v\|^2}{\sigma^2}} \chi \left(\begin{pmatrix} \lambda & w^* \\ 0 & B \end{pmatrix}, \lambda, e_1 \right) |\det(B - \lambda I_{n-1})|^2 \right), \end{aligned} \tag{38}$$

where $v \in \mathbb{P}(\mathbb{C}^n)$ has the uniform distribution, $\hat{y}_v = P_{v^\perp} \hat{A}v / \|v\|$, and $\lambda \sim \mathcal{N}_{\mathbb{C}}(\hat{\lambda}, \sigma^2)$, $w \sim \mathcal{N}_{\mathbb{C}^{n-1}}(\hat{w}, \sigma^2)$, and $B \sim \mathcal{N}_{\mathbb{C}^{(n-1) \times (n-1)}}(\hat{B}, \sigma^2)$ are independent Gaussian random variables centered at

$$\hat{\lambda} := \frac{\langle \hat{A}v, v \rangle}{\|v\|^2}; \quad \hat{w} := J^* U_v^* \hat{A}^* U_v e_1; \quad \hat{B} := J^* U_v^* \hat{A} U_v J.$$

Here, J is the $n \times (n-1)$ matrix whose columns are e_2, \dots, e_n (and, hence, $J^* = (0 \quad \text{Id}_{n-1})$ is the matrix of $P_{e_1^\perp}: \mathbb{C}^n \rightarrow e_1^\perp$). In particular, if $\hat{A} = 0$ then $\hat{y}_v = 0$, $\hat{\lambda} = 0$, $\hat{w} = 0$ and $\hat{B} = 0$, so v can be removed from the expected value in (38).

PROOF. The first claim of the theorem follows directly from Theorem 6.2 and Proposition 6.4.

For the second claim, let I be the left-hand side of (38). Change $\chi(A, \lambda, v)$ to $(\sigma^2 \pi)^{-n^2} \chi(A, \lambda, v) e^{-\frac{\|A - \hat{A}\|_F^2}{\sigma^2}}$ in the first formula to get

$$I = \frac{1}{(\sigma^2 \pi)^{n^2}} \int_{(\lambda, v) \in \mathbb{C} \times \mathbb{P}(\mathbb{C}^n)} \int_{A: Av = \lambda v} \chi(A, \lambda, v) |\det(A_{\lambda, v})|^2 e^{-\frac{\|A - \hat{A}\|_F^2}{\sigma^2}} dA d(\lambda, v).$$

Note that $\{A : Av = \lambda v\}$ can be parametrized by

$$(w, B) \mapsto A = U_v \begin{pmatrix} \lambda & w^* \\ 0 & B \end{pmatrix} U_v^*,$$

where $w \in \mathbb{C}^{n-1}$, $B \in \mathbb{C}^{(n-1) \times (n-1)}$. This parametrization preserves distances; moreover $|\det(A_{\lambda, v})| = |\det(\lambda I_{n-1} - B)|$ and from the fact that χ is unitarily invariant, we have

$$I = \frac{1}{(\sigma^2 \pi)^{n^2}} \int_{(\lambda, v, w, B)} \chi \left(\begin{pmatrix} \lambda & w^* \\ 0 & B \end{pmatrix}, \lambda, e_1 \right) |\det(\lambda I_{n-1} - B)|^2 e^{-\frac{\|A - \hat{A}\|_F^2}{\sigma^2}} d(\lambda, v, w, B),$$

where A is given by the formula above. Note now that we can write

$$\begin{aligned} \|A - \hat{A}\|_F^2 &= \left\| \begin{pmatrix} \lambda & w^* \\ 0 & B \end{pmatrix} - U_v^* \hat{A} U_v \right\|_F^2 \\ &= |\lambda - e_1^* U_v^* \hat{A} U_v e_1|^2 + \|w^* - e_1^* U_v^* \hat{A} U_v J\|^2 + \|B - J^* U_v^* \hat{A} U_v J\|_F^2 \\ &\quad + \|J^* U_v^* \hat{A} U_v e_1\|^2 \\ &= \left| \lambda - \frac{v^* \hat{A} v}{\|v\|^2} \right|^2 + \left\| w - J^* U_v^* \hat{A}^* U_v e_1 \right\|^2 + \|B - J^* U_v^* \hat{A} U_v J\|_F^2 + \|\hat{y}_v\|^2, \end{aligned}$$

and the second claim of the theorem follows noting that the volume of $\mathbb{P}(\mathbb{C}^n)$ is $\pi^{n-1}/\Gamma(n)$. \square

6.4 The linear solution variety

It will be useful to consider a geometrical scheme similar to that of §6.3 for the case of solving linear systems: we consider

$$\mathcal{V}^{\text{lin}} = \{(M, v) \in \mathbb{C}^{(n-1) \times n} \times \mathbb{P}(\mathbb{C}^n) : Mv = 0\}.$$

The linear solution variety \mathcal{V}^{lin} is a $n(n-1)$ -dimensional smooth submanifold of $\mathbb{C}^{(n-1) \times n} \times \mathbb{P}(\mathbb{C}^n)$, and again it inherits the Riemannian structure of the ambient space (cf. [19, (17.14)]).

The linear solution variety is equipped with two projections

$$\begin{aligned} \pi_1^{\text{lin}}: \quad \mathcal{V}^{\text{lin}} &\rightarrow \mathbb{C}^{(n-1) \times n} & \text{and} & \quad \pi_2^{\text{lin}}: \quad \mathcal{V}^{\text{lin}} &\rightarrow \mathbb{P}(\mathbb{C}^n) \\ (M, v) &\mapsto M & (M, v) &\mapsto v. \end{aligned} \quad (39)$$

Note that π_2^{lin} is regular at every $(M, v) \in \mathcal{V}^{\text{lin}}$ and π_1^{lin} is regular at $(M, v) \in \mathcal{V}^{\text{lin}}$ if and only if M is of maximal rank.

For $M \in \mathbb{C}^{(n-1) \times n}$, $(\pi^{\text{lin}})^{-1}(M)$ is a copy of the projective linear subspace corresponding to the kernel of M in $\mathbb{P}(\mathbb{C}^n)$, and for $v \in \mathbb{P}(\mathbb{C}^n)$, $(\pi_2^{\text{lin}})^{-1}(v)$ is a copy of the linear subspace of $\mathbb{C}^{(n-1) \times n}$ consisting of the matrices A such that $Av = 0$.

We can apply Theorem 6.2 for integrating functions in \mathcal{V}^{lin} using the projections in (39). The tangent space to \mathcal{V}^{lin} at a regular point (M, v) can be identified with

$$\{(\dot{M}, \dot{v}) : \dot{M}v + M\dot{v} = 0, v^*\dot{v} = 0\} = \{(\dot{M}, \dot{v}) : \dot{v} = \varphi(\dot{M})\}, \quad \varphi(\dot{M}) = -M^\dagger \dot{M}v.$$

Note that φ is a linear mapping defined from $\mathbb{C}^{(n-1) \times n}$ to v^\perp . A routine computation shows that, if $\|v\| = 1$, then $\varphi\varphi^* : v^\perp \rightarrow v^\perp$ satisfies $\varphi\varphi^*(w) = P_{v^\perp} M^\dagger (M^\dagger)^* i_{\mathbb{C}^n} w$. Writing down the singular value decomposition of M , it follows that $\det(\varphi\varphi^*) = \det(MM^*)^{-1}$. From Lemma 6.3 it follows that

$$\frac{\text{NJ}(\pi^{\text{lin}})(M, v)}{\text{NJ}(\pi_2^{\text{lin}})(M, v)} = |\det(MM^*)|. \quad (40)$$

Proposition 6.6 *Let $\phi^{\text{lin}}: \mathcal{V}^{\text{lin}} \rightarrow [0, \infty]$ be a measurable unitarily invariant function in the sense that $\phi^{\text{lin}}(M, v) = \phi^{\text{lin}}(MU^*, Uv)$ for any unitary matrix $U \in \mathcal{U}_n$. Then,*

$$\mathbb{E}_{M \sim \mathcal{N}_{\mathbb{C}^{(n-1) \times n}}}(\phi^{\text{lin}}(M, \ker(M))) = \frac{1}{\Gamma(n)} \mathbb{E}_{B \sim \mathcal{N}_{\mathbb{C}^{(n-1) \times (n-1)}}}(\phi^{\text{lin}}((0 \ B), e_1) |\det(B)|^2).$$

PROOF. Let $\chi(M) = \phi^{\text{lin}}(M, \ker(M))e^{-\|M\|_F^2}$. Theorem 6.2 and (40) imply that

$$\int_{M \in \mathbb{C}^{(n-1) \times n}} \chi(M) dM = \int_{v \in \mathbb{P}(\mathbb{C}^n)} \int_{M: Mv=0} |\det(MM^*)| \chi(M) dM dv.$$

Now, $|\det(MM^*)|^2 \chi(M) = |\det(MU(MU)^*)|^2 \chi(MU^*)$ for all $U \in \mathcal{U}_n$ by hypothesis. Hence, by parametrizing $\{M : Mv = 0\}$ by $\{(0 \ B)U_v^* : B \in \mathbb{C}^{(n-1) \times (n-1)}\}$ where $U_v \in \mathcal{U}_n$ is any matrix satisfying $U_v e_1 = v$ (we are assuming $\|v\| = 1$), we conclude that the inner integral in the right-hand side of the formula above does not depend on v . We thus have

$$\int_{M \in \mathbb{C}^{(n-1) \times n}} \chi(M) dM = \text{vol}(\mathbb{P}(\mathbb{C}^n)) \int_{B \in \mathbb{C}^{(n-1) \times (n-1)}} |\det(B)|^2 \chi((0, B)) dB.$$

The proposition follows from the form of the Gaussian density (recall §2.6) by noting that $\text{vol}(\mathbb{P}(\mathbb{C}^n)) = \pi^{n-1}/\Gamma(n)$. \square

Assume now that we are given an a.e. continuous function $\alpha: \mathbb{C}^{(n-1) \times (n-1)} \rightarrow [0, \infty]$. We can produce a unitarily invariant function defined on \mathcal{V}^{lin} as follows

$$\phi^{\text{lin}}(M, v) = \mathbb{E}_{Q:(M,Q) \in \mathcal{A}_n} (\alpha(MQ)),$$

where \mathcal{A}_n is given in (22). Note that

$$\phi^{\text{lin}}((0 \ B), e_1) = \mathbb{E}_{U \in \mathcal{U}_{n-1}} (\alpha(BU)).$$

It is a simple exercise to check that ϕ^{lin} is unitarily invariant in the sense of Proposition 6.6. Applying Proposition 6.6 to ϕ^{lin} then yields

$$\mathbb{E}_{M \sim \mathcal{N}_{\mathbb{C}^{(n-1) \times n}}} \mathbb{E}_{Q:(M,Q) \in \mathcal{A}_n} (\alpha(MQ)) = \frac{1}{\Gamma(n)} \mathbb{E}_B \mathbb{E}_{U \in \mathcal{U}_{n-1}} (\alpha(BU) |\det(B)|^2).$$

(The expected value in the unitary group is again with respect to the uniform distribution coming from its Riemannian structure). Finally, using Fubini's theorem, we can interchange the integration order in the right-hand term, and then note that the isometry $B \mapsto BU$ preserves the value of the integral inside. We obtain the following corollary.

Corollary 6.7 *Let $\alpha: \mathbb{C}^{(n-1) \times (n-1)} \rightarrow [0, \infty]$ be an a.e. continuous function. Then,*

$$\mathbb{E}_{M \sim \mathcal{N}_{\mathbb{C}^{(n-1) \times n}}} \mathbb{E}_{Q:(M,Q) \in \mathcal{A}_n} (\alpha(MQ)) = \frac{1}{\Gamma(n)} \mathbb{E}_{B \sim \mathcal{N}_{\mathbb{C}^{(n-1) \times (n-1)}}} (\alpha(B) |\det(B)|^2).$$

□

7 Proof of Theorem 2.14

We begin with the following result.

Proposition 7.1 *The following inequality holds for every $\hat{A} \in \mathbb{C}^{m \times m}$ and $\sigma > 0$.*

$$\mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{m \times m}}(\hat{A}, \sigma^2)} (\|A^{-1}\|_F^2 |\det(A)|^2) \leq \frac{m}{\sigma^2} \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{m \times m}}(\hat{A}, \sigma^2)} (|\det(A)|^2).$$

Furthermore, equality holds if and only if $\hat{A} = 0$. In particular,

$$\mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{m \times m}}(0, \sigma^2)} (\|A^{-1}\|_F^2 |\det(A)|^2) = m! m \sigma^{2m-2}.$$

PROOF. Expanding the determinant of A by the k th column we have

$$\det(A) = \sum_{j=1}^m (-1)^{j+k} a_{j,k} \det A^{j,k},$$

where $A^{j,k}$ denotes the matrix that results from the matrix A by removing the j th row and k th column. Hence,

$$|\det(A)|^2 = \det(A) \overline{\det(A)} = \sum_{j,j'=1}^m (-1)^{j+j'+2k} a_{j,k} \overline{a_{j',k}} \det A^{j,k} \overline{\det A^{j',k}}.$$

Observe that the random variables $a_{j,k}$ and $a_{j',k}$ are independent of $\det A^{j,k}$ and $\det A^{j',k}$. Then,

$$\mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{m \times m}}(\hat{A}, \sigma^2)} |\det(A)|^2 = \sum_{j,j'=1}^m (-1)^{j+j'+2k} \mathbb{E}(a_{j,k} \overline{a_{j',k}}) \mathbb{E}(\det A^{j,k} \overline{\det A^{j',k}}).$$

Now observe that

$$\mathbb{E}(a_{j,k} \overline{a_{j',k}}) = \begin{cases} \hat{a}_{j,k} \overline{\hat{a}_{j',k}} & \text{if } j \neq j'; \\ \sigma^2 + |\hat{a}_{j,k}|^2 & \text{otherwise.} \end{cases}$$

We conclude that for $k = 1, \dots, m$,

$$\mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{m \times m}}(\hat{A}, \sigma^2)} |\det(A)|^2 = \mathbb{E} |\det([A; k; \hat{A}_k])|^2 + \sigma^2 \sum_{j=1}^m \mathbb{E} |\det A^{j,k}|^2, \quad (41)$$

where $[A; k; \hat{A}_k]$ is the matrix formed by replacing the (random) k th column of A by the (deterministic) k th column of \hat{A} . Summing on k we get

$$m \mathbb{E}_A |\det(A)|^2 = \sum_{k=1}^m \mathbb{E} |\det([A; k; \hat{A}_k])|^2 + \sigma^2 \sum_{j,k=1}^m \mathbb{E} |\det A^{j,k}|^2. \quad (42)$$

On the other hand, from a direct application of Cramer's Rule and (42), we deduce that

$$\begin{aligned} \sigma^2 \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{m \times m}}(\hat{A}, \sigma^2)} \|A^{-1}\|_F^2 |\det(A)|^2 &= \sigma^2 \sum_{j,k=1}^m \mathbb{E} |\det A^{j,k}|^2 \\ &= m \mathbb{E} |\det(A)|^2 - \sum_{k=1}^m \mathbb{E} |\det([A; k; \hat{A}_k])|^2, \end{aligned}$$

and the first claim of the proposition follows. Moreover, when $\hat{A} = 0$, the last term in the sum above is zero. We leave the proof of the converse to the reader. Using (41) and the fact that the matrices $A^{j,k}$ are $\mathcal{N}_{\mathbb{C}^{(m-1) \times (m-1)}}(0, \sigma^2)$ -distributed, one can prove working by induction the equality

$$\mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{m \times m}}(0, \sigma^2)} |\det(A)|^2 = \sigma^{2m} m!.$$

The second claim of the proposition follows. \square

Corollary 7.2

$$\mathbb{E}_{M \sim \mathcal{N}_{\mathbb{C}^{(n-1) \times n}}} (\|M^\dagger\|_F^2) = n - 1.$$

PROOF. From Proposition 6.6 with $\phi^{\text{lin}}(M, \zeta) = \|M^\dagger\|_F^2$, we have

$$\mathbb{E}_{M \sim \mathcal{N}_{\mathbb{C}^{(n-1) \times n}}} (\|M^\dagger\|_F^2) = \frac{1}{\Gamma(n)} \mathbb{E}_{B \sim \mathcal{N}_{\mathbb{C}^{(n-1) \times (n-1)}}} (\|B^{-1}\|_F^2 |\det(B)|^2) \xrightarrow{\text{Prop. 7.1}} n - 1$$

as claimed. \square

Corollary 7.3 For any $\hat{B} \in \mathbb{C}^{(n-1) \times (n-1)}$, $\sigma > 0$, and $\lambda \in \mathbb{C}$, we have

$$\mathbb{E}_B (\|(B - \lambda I_{n-1})^{-1}\|_F^2 |\det(B - \lambda I_{n-1})|^2) \leq \frac{n-1}{\sigma^2} \mathbb{E}_B (|\det(B - \lambda I_{n-1})|^2),$$

where $B \sim \mathcal{N}_{\mathbb{C}^{(n-1) \times (n-1)}}(\hat{B}, \sigma^2)$.

PROOF. Note that

$$\mathbb{E}_B (\|(B - \lambda I_{n-1})^{-1}\|_F^2 |\det(B - \lambda I_{n-1})|^2) = \mathbb{E}_{C \sim \mathcal{N}_{\mathbb{C}^{(n-1) \times (n-1)}}(\hat{C}, \sigma^2)} (\|C^{-1}\|_F^2 |\det C|^2),$$

where $\hat{C} = \hat{B} - \lambda I_{n-1}$. The proof readily follows from Proposition 7.1. \square

PROOF OF THEOREM 2.14. Fix any a.e. continuous mapping $v \mapsto U_v$ such that for $v \in \mathbb{P}(\mathbb{C}^n)$, U_v is a unitary matrix with $U_v e_0 = v/\|v\|$. From (38) applied to $\chi(A, \lambda, v) = \frac{1}{n} \mu_F^2(A, \lambda, v) / \|A\|_F^2 = \frac{1}{n} \|A_{\lambda, v}^{-1}\|_F^2$ we have

$$\begin{aligned} & \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}(\hat{A}, \sigma^2)} \left(\frac{\mu_{F, \mathbf{av}}^2(A)}{\|A\|_F^2} \right) \\ &= \frac{1}{n \Gamma(n) \sigma^{2(n-1)}} \mathbb{E}_v \mathbb{E}_{(\lambda, w, B)} (e^{-\frac{\|\hat{y}_v\|^2}{\sigma^2}} \|(B - \lambda I_{n-1})^{-1}\|_F^2 |\det(B - \lambda I_{n-1})|^2), \end{aligned} \tag{43}$$

where $\hat{y}_v = P_{v^\perp} \hat{A} v / \|v\|$, $v \in \mathbb{P}(\mathbb{C}^n)$ has the uniform distribution, and $\lambda \sim \mathcal{N}_{\mathbb{C}}(\hat{\lambda}, \sigma^2)$, $w \sim \mathcal{N}_{\mathbb{C}^{n-1}}(\hat{w}, \sigma^2)$, and $B \sim \mathcal{N}_{\mathbb{C}^{(n-1) \times (n-1)}}(\hat{B}, \sigma^2)$ for some $\hat{\lambda}, \hat{w}, \hat{B}$ which depend uniquely on \hat{A} and v .

From (43) and Corollary 7.3 we have

$$\begin{aligned} & \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}(\hat{A}, \sigma^2)} \left(\frac{\mu_{F, \mathbf{av}}^2(A)}{\|A\|_F^2} \right) \\ & \leq \frac{n-1}{\sigma^2 \sigma^{2(n-1)} \Gamma(n+1)} \mathbb{E}_v \mathbb{E}_{(\lambda, w, B)} (e^{-\frac{\|\hat{y}_v\|^2}{\sigma^2}} |\det(B - \lambda I_{n-1})|^2). \end{aligned}$$

Now, if we apply again (38) to the constant function $\chi \equiv 1/n$ we get

$$1 = \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{m \times m}}(\hat{A}, \sigma^2)} (1) = \frac{1}{\Gamma(n+1)\sigma^{2(n-1)}} \mathbb{E}_v \mathbb{E}_{(\lambda, w, B)} (e^{-\frac{\|\hat{y}_v\|^2}{\sigma^2}} |\det(B - \lambda I_{n-1})|^2),$$

and we have thus proved that

$$\mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}(\hat{A}, \sigma^2)} \left(\frac{\mu_{F, \text{av}}^2(A)}{\|A\|_F^2} \right) \leq \frac{n-1}{\sigma^2} \leq \frac{n}{\sigma^2},$$

as claimed.

This proves the first part of Theorem 2.14. For the second part of the theorem, let

$$I = \mathbb{E}_{A \sim \mathcal{U}(\mathbb{S})} \mu_{F, \text{av}}^2(A)$$

be the quantity we want to compute. From the first part of the theorem (with $\hat{A} = 0$ and $\sigma = 1$) we have

$$\frac{1}{\pi^{n^2}} \int_{A \in \mathbb{C}^{n \times n}} \frac{\mu_{F, \text{av}}^2(A)}{\|A\|_F^2} e^{-\|A\|_F^2} dA \leq n. \quad (44)$$

On the other hand,

$$\frac{1}{\pi^{n^2}} \int_{A \in \mathbb{C}^{n \times n}} \frac{\mu_{F, \text{av}}^2(A)}{\|A\|_F^2} e^{-\|A\|_F^2} dA = \frac{1}{\pi^{n^2}} \int_0^\infty \frac{e^{-\rho^2}}{\rho^2} \int_{A: \|A\|_F = \rho} \mu_{F, \text{av}}^2(A) dA d\rho. \quad (45)$$

Now, because $\mu_{F, \text{av}}(A)$ is invariant under multiplication of A by nonzero complex numbers, denoting $\nu_\rho = \text{vol}(A : \|A\|_F = \rho)$, we have

$$\frac{1}{\nu_\rho} \int_{A: \|A\|_F = \rho} \mu_{F, \text{av}}^2(A) dA = I, \quad 0 < \rho < \infty. \quad (46)$$

We deduce from (44–46) that

$$\frac{I}{\pi^{n^2}} \int_0^\infty \frac{\nu_\rho e^{-\rho^2}}{\rho^2} d\rho \leq n.$$

Note now that

$$\nu_\rho = \frac{2\pi^{n^2}}{\Gamma(n^2)} \rho^{2n^2-1}$$

to conclude that

$$I \leq \frac{n\Gamma(n^2)}{2 \int_0^\infty \rho^{2n^2-3} e^{-\rho^2} d\rho} = \frac{n\Gamma(n^2)}{\Gamma(n^2-1)} = n(n^2-1) \leq n^3.$$

The theorem follows. \square

8 Proof of Propositions 2.22 and 2.24

8.1 Proof of Proposition 2.22

First note that Path-follow starts by normalizing the input, so from (14) we can assume that $\|A_0\|_F = 1$ and $A \sim \mathcal{N}_{\mathbb{C}^{n \times n}, T}(0, 1)$ where $T = \sqrt{2}n$. From Remark 2.23, for integration purposes we can also assume that $(\mathcal{L}_{A_0, A} \setminus \{A_0\}) \cap \Sigma = \emptyset$. Corollary 2.21 with $q = 1$ implies that

$$\begin{aligned} \text{Avg_Num_Iter}(A_0, \lambda_0, v_0) &= \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}, T}} [K(A, A_0, \lambda_0, v_0)] \\ &\leq 2 + \sqrt{6}C \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}, T}} \|A\|_F \int_{t_1}^1 \frac{\mu^2(A_t, \lambda_t, v_t)}{\|A_t\|_F^2} dt \quad (47) \\ &\leq 2 + \sqrt{6}C \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}, T}} \|A\|_F \int_{t_1}^1 \sum_{j=1}^n \frac{\mu^2(A_t, \lambda_t^{(j)}, v_t^{(j)})}{\|A_t\|_F^2} dt, \end{aligned}$$

where $A_t = (1-t)A_0 + tA$, the pairs $(\lambda_t^{(j)}, v_t^{(j)})$ are defined by continuation for all the eigenpairs of A_t , and

$$t_1 = \inf_{\|A\|_F \leq T} \frac{1}{\|A\|_F (\sin \alpha \cot(\text{Choose_step}(A_0, \dot{A}_0, \lambda_0, v_0)) - \cos \alpha) + 1}.$$

We therefore have (use $T = \sqrt{2}n$)

$$\begin{aligned} \text{Avg_Num_Iter}(A_0, \lambda_0, v_0) &\leq 2 + \sqrt{6}CnT \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}, T}} \left(\int_{t_1}^1 \frac{\mu_{\text{av}}^2(A_t)}{\|A_t\|_F^2} dt \right) \\ &\leq 2 + \sqrt{48}Cn^2 \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}} \left(\int_{t_1}^1 \frac{\mu_{\text{av}}^2(A_t)}{\|A_t\|_F^2} dt \right). \quad (48) \end{aligned}$$

We have used (13) and Lemma 2.16 for the last inequality. In order to bound the last term in the previous expression, we interchange the order of integration,

$$\mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}} \left(\int_{t_1}^1 \frac{\mu_{\text{av}}^2(A_t)}{\|A_t\|_F^2} dt \right) = \int_{t_1}^1 \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}} \left(\frac{\mu_{\text{av}}^2(A_t)}{\|A_t\|_F^2} \right) dt.$$

Now, for fixed t , if $A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}$ then $A_t = (1-t)A_0 + tA$ satisfies $A_t \sim \mathcal{N}_{\mathbb{C}^{n \times n}}((1-t)A_0, t^2)$ and from Theorem 2.14 we have

$$\mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}} \left(\frac{\mu_{\text{av}}^2(A_t)}{\|A_t\|_F^2} \right) \leq \frac{n}{t^2},$$

which implies

$$\mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}} \left(\int_{t_1}^1 \frac{\mu_{\text{av}}^2(A_t)}{\|A_t\|_F^2} dt \right) \leq \int_{t_1}^1 \frac{n}{t^2} dt \leq \frac{n}{t_1}. \quad (49)$$

We are thus left with the task of evaluating t_1 . Note that $s_1 = \text{Choose_step}(A_0, \dot{A}_0, \lambda_0, v_0)$ (the length of the first step in the execution of **Path-follow**) is at least $\frac{R}{\mu^2(A_0, \lambda_0, v_0)}$ by Proposition 5.8. Hence $\cot s_1 \leq \frac{1}{s_1} \leq \frac{\mu^2(A_0, \lambda_0, v_0)}{R}$ and it follows that

$$t_1 \geq \frac{1}{T(\cot s_1 + 1) + 1} \geq \Omega\left(\frac{1}{n\mu^2(A_0, \lambda_0, v_0)}\right).$$

Putting together this bound and inequalities (48) and (49) we deduce the claimed bound for $\text{Avg_Num_Iter}(A_0, \lambda_0, v_0)$.

We next prove the smoothed analysis bounds. Reasoning as in (47) we see that the smoothed number of iterations $\text{Smd_Num_Iter}(A_0, \lambda_0, v_0, \sigma)$ is bounded by

$$2 + \sqrt{6}C \sup_{\hat{A} \in \mathbb{S}} \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}, T}(\hat{A}, \sigma^2)} \sum_{j=1}^n \|A\|_F \int_{t_1}^1 \frac{\mu^2(A_t, \lambda_t^{(j)}, v_t^{(j)})}{\|A_t\|_F^2} dt.$$

The rest of the argument is almost exactly as above, the only difference being the bound $\|A\|_F \leq T + \|\hat{A}\|_F = \sqrt{2}n + 1$. \square

8.2 Proof of Proposition 2.24

We are now following all the n paths (each starting with a different eigenpair of A_0). Applying Corollary 2.21 with $q = 1$ to each of them we obtain

$$\begin{aligned} \text{Avg_Num_Iter}(A_0) &= \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}, T}} \sum_{j=1}^n K(A, A_0, \lambda^{(j)}, v^{(j)}) \\ &\leq 2n + \sqrt{6}C \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}, T}} \|A\|_F \sum_{j=1}^n \int_{t_1^{(j)}}^1 \frac{\mu^2(A_t, \lambda_t^{(j)}, v_t^{(j)})}{\|A_t\|_F^2} dt \quad (50) \\ &\leq 2n + \sqrt{6}C \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}, T}} \|A\|_F \int_{t_1^*}^1 \sum_{j=1}^n \frac{\mu^2(A_t, \lambda_t^{(j)}, v_t^{(j)})}{\|A_t\|_F^2} dt, \end{aligned}$$

where $t_1^* = \min\{t_1^{(1)}, \dots, t_1^{(n)}\}$. We can now reason as in the preceding proof to deduce that

$$\text{Avg_Num_Iter}(A_0) = \mathcal{O}\left(\frac{n^3}{t_1^*}\right)$$

as well as the fact that, for $j = 1, \dots, n$, $t_1^{(j)} \geq \Omega\left(\frac{1}{n\mu^2(A_0, \lambda^{(j)}, v^{(j)})}\right)$. It follows from these bounds that

$$t_1^* \geq \Omega\left(\frac{1}{n\mu_{\max}^2(A_0)}\right).$$

The rest of the proof follows as in the preceding proposition. \square

9 Proof of Theorem 2.29

We begin with an auxiliary result. For simplicity, in what follows we write $\mathbb{S} := \mathbb{S}(\mathbb{C}^{n \times n})$. We also consider the manifold

$$\mathcal{S} = \{(A, \dot{A}) \in \mathbb{S} \times \mathbb{S} : \dot{A} \in T_A \mathbb{S}\}$$

and denote by $\mathcal{U}(\mathcal{S})$ the normalized product distribution on it, that is, the probability distribution coming from the product structure. Given any measurable mapping $\phi: \mathcal{S} \rightarrow [0, \infty]$, let

$$I_\phi := \mathbb{E}_{A_0, A \in \mathbb{S}} \left(\int_0^{d_{\mathbb{S}}(A_0, A)} \phi(A_s, \dot{A}_s) ds \right), \quad (51)$$

where, as usual, the A_s are such that $\{A_s : 0 \leq s \leq d_{\mathbb{S}}(A_0, A)\} = \mathcal{L}_{A_0, A}$, and \dot{A}_s is the unit tangent vector (in the direction of the parametrization) to $\mathcal{L}_{A_0, A}$ at A_s .

Lemma 9.1 *For any measurable mapping $\phi: \mathbb{S} \times \mathbb{S} \rightarrow [0, \infty]$ we have*

$$I_\phi = \frac{\pi}{2} \mathbb{E}_{(A, \dot{A}) \sim \mathcal{U}(\mathcal{S})} \phi(A, \dot{A}). \quad (52)$$

PROOF. We consider the manifold $\mathcal{R} = \{(A_0, A, s) \in \mathbb{S} \times \mathbb{S} \times (0, \pi) : s < d_{\mathbb{S}}(A_0, A)\}$ with the product structure and let

$$\begin{aligned} \psi: \quad \mathcal{R} &\rightarrow \mathcal{S} \\ (A_0, A, s) &\mapsto (A_s, \dot{A}_s). \end{aligned}$$

We can then write $I_\phi = \text{vol}(\mathbb{S})^{-2} \int_{\mathcal{R}} \phi \circ \psi$. Applying the coarea formula (Theorem 6.1) yields

$$I_\phi = \frac{1}{\text{vol}(\mathbb{S})^2} \int_{(A, \dot{A}) \in \mathcal{S}} \phi(A, \dot{A}) q(A, \dot{A}) d(A, \dot{A}), \quad (53)$$

where

$$q(A, \dot{A}) = \int_{(B_0, B, s) \in \psi^{-1}(A, \dot{A})} \frac{1}{\text{NJ}(\psi)(B_0, B, s)} d(B_0, B, s).$$

Our goal now is to prove that $q(A, \dot{A})$ is a constant (independent of A, \dot{A}). It shall be useful to consider the two diagonal matrices $\Delta = E_{11}$ and $\dot{\Delta} = E_{22}$ where E_{ij} denotes the standard basis of $\mathbb{C}^{n \times n}$. Note that $(\Delta, \dot{\Delta}) \in \mathcal{S}$.

We now fix $(A, \dot{A}) \in \mathcal{S}$. Let $\sigma: \mathbb{C}^{n \times n} \rightarrow \mathbb{C}^{n \times n}$ be an isometric change of basis such that $\sigma(A) = \Delta$ and $\sigma(\dot{A}) = \dot{\Delta}$. Denoting $\sigma_{\mathcal{S}} = \sigma \times \sigma$ and $\sigma_{\mathcal{R}} = \sigma \times \sigma \times \text{Id}_{\mathbb{R}}$, where $\text{Id}_{\mathbb{R}}$ is the identity mapping in \mathbb{R} , it is easy to check by writing down the formula for A_s that $\psi \circ \sigma_{\mathcal{R}} = \sigma_{\mathcal{S}} \circ \psi$. We are under the hypothesis of [14, Lemma 4,

p. 244] (which holds for surjective maps in general, not only for projections), proving that

$$\text{NJ}(\psi)(\sigma_{\mathcal{R}}^{-1}(C_0, C, s)) = \text{NJ}(\psi)(C_0, C, s), \quad \forall (C_0, C, s) \in \mathcal{R}. \quad (54)$$

Moreover, the mapping $\sigma_{\mathcal{R}}|_{\psi^{-1}(A, \dot{A})}$ is an isometry from $\psi^{-1}(A, \dot{A})$ to $\psi^{-1}(\Delta, \dot{\Delta})$, which from the change of variables theorem implies

$$\begin{aligned} q(A, \dot{A}) &= \int_{(B_0, B, s) \in \psi^{-1}(A, \dot{A})} \frac{1}{\text{NJ}(\psi)(B_0, B, s)} d(B_0, B, s) \\ &\stackrel{(54)}{=} \int_{(B_0, B, s) \in \psi^{-1}(A, \dot{A})} \frac{1}{\text{NJ}(\psi)(\sigma_{\mathcal{R}}^{-1}(B_0, B, s))} d(B_0, B, s) \\ &= \int_{(B_0, B, s) \in \psi^{-1}(\Delta, \dot{\Delta})} \frac{1}{\text{NJ}(\psi)(B_0, B, s)} d(B_0, B, s) = q(\Delta, \dot{\Delta}), \end{aligned}$$

which proves that $q(A, \dot{A})$ is equal to some constant \hat{C} .

Since this holds for any measurable function ϕ , we can take the latter to be constant with value 1 to derive the value of \hat{C} . Then (53) becomes

$$I_{\phi} = \frac{1}{\text{vol}(\mathbb{S})^2} \int_{(A, \dot{A}) \in \mathcal{S}} \hat{C} d(A, \dot{A}) = \frac{\text{vol}(\mathcal{S})}{\text{vol}(\mathbb{S})^2} \hat{C}.$$

And it follows from (51) (always with $\phi \equiv 1$) that

$$I_{\phi} = \mathbb{E}_{A, A_0 \in \mathbb{S}} d_{\mathbb{S}}(A_0, A).$$

Hence,

$$\mathbb{E}_{A, A_0 \in \mathbb{S}} d_{\mathbb{S}}(A, A_0) = \hat{C} \frac{\text{vol}(\mathcal{S})}{\text{vol}(\mathbb{S})^2}.$$

Note that the change of variables $A_0 \mapsto -A_0$ does not change the expected value in this last formula. Moreover, $d_{\mathbb{S}}(A, A_0) + d_{\mathbb{S}}(A, -A_0) = \pi$ for all $A, A_0 \in \mathbb{S}$. Thus,

$$\begin{aligned} \frac{\text{vol}(\mathcal{S})}{\text{vol}(\mathbb{S})^2} 2\hat{C} &= \mathbb{E}_{A, A_0 \in \mathbb{S}} d_{\mathbb{S}}(A, A_0) + \mathbb{E}_{A, A_0 \in \mathbb{S}} d_{\mathbb{S}}(A, -A_0) \\ &= \mathbb{E}_{A, A_0 \in \mathbb{S}} (d_{\mathbb{S}}(A, A_0) + d_{\mathbb{S}}(A, -A_0)) = \pi. \end{aligned}$$

proving that

$$\hat{C} = \frac{\text{vol}(\mathbb{S})^2 \pi}{2\text{vol}(\mathcal{S})},$$

From (53) we conclude that for any measurable nonnegative function ϕ we have

$$I_{\phi} = \frac{\pi}{2} \mathbb{E}_{(A, \dot{A}) \sim \mathcal{U}(\mathcal{S})} \phi(A, \dot{A}),$$

as wanted. □

PROOF OF THEOREM 2.29. Consider the measurable function $\phi: \mathcal{S} \rightarrow [0, \infty]$ defined by

$$\phi(A, \dot{A}) = \frac{1}{n} \sum_{(\lambda, v): Av = \lambda v} \mu(A, \lambda, v) \|(\dot{A}, \dot{\lambda}, \dot{v})\|,$$

for $A \in \mathbb{S}$ and $\dot{A} \in T_A \mathbb{S}$ such that $A \notin \Sigma$, where $\dot{\lambda}, \dot{v}$ are the functions of (A, \dot{A}) and (λ, v) given in Lemma 3.2. (If $A \in \Sigma$ we set $\phi(A, \dot{A}) = \infty$.)

From Theorem 2.19, denoting $I = I_\phi$, we have (for some constant $C > 1$) that

$$\mathbb{E}_{A, A_0 \sim \mathcal{N}_{\mathbb{C}^{n \times n}}} \left(\frac{1}{n} \sum_{\lambda_0, v_0: A_0 v_0 = \lambda_0 v_0} K(A, A_0, \lambda_0, v_0) \right) \leq C I. \quad (55)$$

Note that the left-hand side of (55) is the quantity to be bounded in Theorem 2.29. It is therefore enough for us to show that $I \leq 4n^2$. To do so, write $\mathcal{S}_A := \{A' : (A, A') \in \mathcal{S}\} \subseteq T_A \mathbb{S}$, for $A \in \mathbb{S}$. First note that \mathcal{S}_A is just the unit sphere in $T_A \mathbb{S}$, so it has a natural volume form inherited from $\mathbb{C}^{n \times n}$ and $\text{vol}(\mathcal{S}_A)$ is independent of A . Moreover, the Normal Jacobian of the projection $\mathcal{S} \rightarrow \mathbb{S}$, $(A, \dot{A}) \mapsto A$, is constant and equal to $1/\sqrt{2}$ (this is easy to prove: check that for $\dot{A} \in T_A \mathbb{S}$ the pair of the form (\dot{A}, \dot{A}') in $T_{(A, \dot{A}')} \mathcal{S}$ which is orthogonal to the kernel of the derivative of the projection is $(\dot{A}, -\mathbb{R}e\langle A', \dot{A} \rangle A)$, then note that the vectors of that form obtained from any o.g. basin of $T_A \mathbb{S}$ whose first element is A' are orthogonal and only one of them, $(A', -A)$, changes its norm by $\sqrt{2}$), so we have $\text{vol}(\mathcal{S}) = \sqrt{2}\text{vol}(\mathbb{S})\text{vol}(\mathcal{S}_A)$. From Lemma 9.1 and Theorem 6.1, we then have

$$\begin{aligned} I &= \frac{\pi}{2} \mathbb{E}_{(A, \dot{A}) \sim \mathcal{U}(\mathcal{S})} \phi(A, \dot{A}) \\ &= \frac{\pi}{\sqrt{2}} \mathbb{E}_{A \in \mathbb{S}} \left(\frac{1}{n} \sum_{(\lambda, v): Av = \lambda v} \mu(A, \lambda, v) \mathbb{E}_{\dot{A} \in \mathcal{S}_A} (\|(\dot{A}, \dot{\lambda}, \dot{v})\|) \right). \end{aligned} \quad (56)$$

In order to estimate this last quantity we shall use Lemma 9.2 below. Note first that from Cauchy-Schwartz,

$$\begin{aligned} \mathbb{E}_{\dot{A} \in \mathcal{S}_A} (\|(\dot{A}, \dot{\lambda}, \dot{v})\|) &\leq \left(1 + \mathbb{E}_{\dot{A} \in \mathcal{S}_A} (|\dot{\lambda}|^2 + \|\dot{v}\|^2) \right)^{1/2} \\ &\stackrel{\text{Lemma 9.2}}{\leq} \left(1 + \frac{1}{n^2 - \frac{1}{2}} (1 + 2\mu_F(A, \lambda, v)^2) \right)^{1/2} \\ &\stackrel{n \geq 2}{\leq} \frac{1}{\sqrt{7}} \left(9 + \frac{16}{n^2} \mu_F(A, \lambda, v)^2 \right)^{1/2}. \end{aligned}$$

It is a simple exercise to check that for positive $x \in \mathbb{R}$ we have $x(9 + 16x^2/n^2)^{1/2} \leq 9n/8 + 4x^2/n$. Using this inequality we get from the equations above:

$$I \leq \frac{\pi}{\sqrt{14}} \mathbb{E} \left(\frac{1}{n} \sum_{(\lambda, v): Av = \lambda v} \left(\frac{9n}{8} + \frac{4}{n} \mu_F(A, \lambda, v)^2 \right) \right).$$

We next use Theorem 2.14 (averaging over $A \in \mathbb{S}$) and bound this last quantity by

$$I \leq \frac{\pi}{\sqrt{14}} \left(\frac{9n}{8} + \frac{4}{n} n^3 \right) = \frac{\pi}{\sqrt{14}} \left(\frac{9n}{8} + 4n^2 \right) \underset{n \geq 2}{\leq} 4n^2, \quad (57)$$

which finishes the proof. \square

We have used the following technical lemma which is in the spirit of [2]. Note that as pointed out in the proof of Theorem 2.29, \mathcal{S}_A is just the unit sphere in the tangent space $T_A \mathbb{S}$ and thus has a natural measure inherited from the inner product in $\mathbb{C}^{n \times n}$.

Lemma 9.2 *Let $(A, \lambda, v) \in \mathcal{V}$, $\|A\|_F = 1$. Define $\mathcal{S}_A := \{A' : (A, A') \in \mathcal{S}\} \subseteq T_A \mathbb{S}$ as in the proof of Theorem 2.29 (that is, \mathcal{S}_A is the unit sphere of $T_A \mathbb{S}$) and, for $\dot{A} \in \mathcal{S}_A$, let $\dot{\lambda}, \dot{v}$ be as in Lemma 3.2. Then,*

$$\mathbb{E}_{\dot{A} \in \mathcal{S}_A} (|\dot{\lambda}|^2) = \frac{1}{n^2 - \frac{1}{2}} \left(\mu_\lambda(A, \lambda, v)^2 - \frac{|\lambda|^2}{2} \right),$$

and

$$\mathbb{E}_{\dot{A} \in \mathcal{S}_A} (\|\dot{v}\|^2) = \frac{1}{n^2 - \frac{1}{2}} \|A_{\lambda, v}^{-1}\|_F^2.$$

In particular, from Proposition 3.3,

$$\mathbb{E}_{\dot{A} \in \mathcal{S}_A} (|\dot{\lambda}|^2 + \|\dot{v}\|^2) \leq \frac{1}{n^2 - \frac{1}{2}} \left(1 + 2\|A_{\lambda, v}^{-1}\|_F^2 \right) = \frac{1}{n^2 - \frac{1}{2}} (1 + 2\mu_F(A, \lambda, v)^2).$$

PROOF. Note that $T_A \mathbb{S}$ coincides with the (real) orthogonal complement, with respect to $\Re \langle \cdot, \cdot \rangle_F$, of $A \in \mathbb{C}^{n \times n}$. Thus $\dim_{\mathbb{R}}(T_A \mathbb{S}) = 2n^2 - 1$. On this space we consider the push-forward measure of the standard Gaussian distribution on $\mathbb{C}^{n \times n}$ by the orthogonal projection $\mathbb{C}^{n \times n} \rightarrow T_A \mathbb{S}$.

Since $T_A \mathbb{S}$ may be split in the (real) orthogonal decomposition $\mathbb{R}\sqrt{-1}A \oplus A^\perp$, where $\mathbb{R}\sqrt{-1}A$ is the linear real subspace generated by $\sqrt{-1}A$, in particular we conclude that the Gaussian distribution on $T_A \mathbb{S}$ coincides with the distribution $t\sqrt{-1}A + \dot{B} \in T_A \mathbb{S}$ where $t \sim \mathcal{N}(0, \frac{1}{2})$ and $\dot{B} \sim \mathcal{N}_{A^\perp}$ are independent.

Claim I: Given a linear operator $L : T_A \mathbb{S} \rightarrow \mathbb{C}^k$, we have

$$\mathbb{E}_{\dot{A} \in T_A \mathbb{S}} (\|L(\dot{A})\|^2) = \left(n^2 - \frac{1}{2} \right) \mathbb{E}_{\dot{A} \in \mathcal{S}_A} (\|L(\dot{A})\|^2);$$

The claim follows integrating in polar coordinates. More precisely,

$$\begin{aligned}
\mathbb{E}_{\dot{A} \in T_A \mathbb{S}} (\|L(\dot{A})\|^2) &= \frac{1}{\pi^{n^2 - \frac{1}{2}}} \int_{\dot{A} \in T_A \mathbb{S}} \|L(\dot{A})\|^2 e^{-\|\dot{A}\|^2} d\dot{A} \\
&= \frac{1}{\pi^{n^2 - \frac{1}{2}}} \int_0^{+\infty} \rho^{2n^2} e^{-\rho^2} d\rho \cdot \int_{\dot{A} \in \mathcal{S}_A} \|L(\dot{A})\|^2 d\dot{A} \\
&= \left(n^2 - \frac{1}{2} \right) \mathbb{E}_{\dot{A} \in \mathcal{S}_A} (\|L(\dot{A})\|^2),
\end{aligned}$$

where we have used that $\int_0^{+\infty} \rho^{2n^2} e^{-\rho^2} d\rho = \frac{1}{2} \Gamma(n^2 + \frac{1}{2})$, and $\text{vol}(\mathcal{S}_A) = 2\pi^{n^2 - \frac{1}{2}} / \Gamma(n^2 - \frac{1}{2})$.

Claim II: For $\|v\| = 1$, the push-forward measure of the Gaussian distribution on A^\perp by the map $f: A^\perp \rightarrow v^\perp$, $\dot{A} \mapsto P_{v^\perp}(\dot{A}v)$, is the standard Gaussian on v^\perp .

Note that for all $B \in \mathbb{C}^{n \times n}$, we have $\langle uv^*, B \rangle_F = \text{tr}(B^* uv^*) = v^* B^* u = \langle u, Bv \rangle$. Then, the set $F = \{\dot{w}v^* : \dot{w} \in v^\perp\}$ is a linear subspace of A^\perp , and the kernel of f is the Hermitian complement of F as subset of A^\perp . Since $f|_F: F \rightarrow v^\perp$ is a linear isometry, the claim follows from the characterization of the standard Gaussian distribution.

Claim III: Let $m \in \mathbb{N}$. If $\eta \sim \mathcal{N}_{\mathbb{C}^m}$, and $x \in \mathbb{C}^m$ then

$$\mathbb{E}_{\eta \sim \mathcal{N}_{\mathbb{C}^m}} (|\langle \eta, x \rangle|^2) = \|x\|^2.$$

The proof of this claim is a standard exercise and is left to the reader.

Now we are ready to prove the lemma. We chose a representative of v such that $\|v\| = 1$ and a representative of the left eigenvector u such that $\langle u, v \rangle = 1$. Note that this implies by Proposition 3.3 and Lemma 3.2:

$$\mu_\lambda(A, \lambda, v) = \frac{\|u\| \|v\|}{|\langle u, v \rangle|} = \|u\|, \quad \dot{\lambda} = \langle \dot{A}v, u \rangle, \quad \dot{v} = -A_{\lambda, v}^{-1} P_{v^\perp} \dot{A}v.$$

For the first statement, we have

$$\begin{aligned}
\mathbb{E}_{\dot{A} \in T_A \mathbb{S}} (|\dot{\lambda}|^2) &= \mathbb{E}_{\dot{A} \in T_A \mathbb{S}} (|\langle \dot{A}v, u \rangle|^2) = \mathbb{E}_{\dot{A} \in T_A \mathbb{S}} (|\langle \dot{A}, uv^* \rangle_F|^2) \\
&= \mathbb{E}_{\dot{B} \in \mathcal{N}_{A^\perp}} \mathbb{E}_{t \sim \mathcal{N}(0, \frac{1}{2})} (|\langle t\sqrt{-1}A + \dot{B}, uv^* \rangle_F|^2).
\end{aligned}$$

Since the mixed term of the expansion of $|\langle t\sqrt{-1}A + \dot{B}, uv^* \rangle_F|^2$ is linear in t , its expected value is zero. Hence,

$$\begin{aligned}
\mathbb{E}_{\dot{A} \in T_A \mathbb{S}} (|\langle \dot{A}v, u \rangle|^2) &= \mathbb{E}_{\dot{B} \in \mathcal{N}_{A^\perp}} \mathbb{E}_{t \sim \mathcal{N}(0, \frac{1}{2})} (t^2 |\langle A, uv^* \rangle_F|^2 + |\langle \dot{B}, uv^* \rangle_F|^2) \\
&= \frac{|\lambda|^2}{2} + \mathbb{E}_{\dot{B} \in \mathcal{N}_{A^\perp}} (|\langle \dot{B}, \pi_{A^\perp}(uv^*) \rangle_F|^2).
\end{aligned}$$

where we have denoted by $\pi_{A^\perp}(uv^*)$ the orthogonal projection of uv^* onto A^\perp . With the identification of A^\perp and \mathbb{C}^{n^2-1} as Hermitian spaces, from Claim III (with $m = n^2 - 1$) we conclude that

$$\mathbb{E}_{\dot{B} \in \mathcal{N}_{A^\perp}} (|\langle \dot{B}, \pi_{A^\perp}(uv^*) \rangle_F|^2) = \|\pi_{A^\perp}(uv^*)\|_F^2 = \|uv^*\|_F^2 - |\lambda|^2 = \|u\|^2 - |\lambda|^2.$$

We have then proved:

$$\mathbb{E}_{\dot{A} \in T_A \mathbb{S}} (|\langle \dot{A}v, u \rangle|^2) = \mu_\lambda(A, \lambda, v)^2 - \frac{|\lambda|^2}{2},$$

and from claim I we conclude:

$$\mathbb{E}_{\dot{A} \in \mathcal{S}_A} (|\langle \dot{A}v, u \rangle|^2) = \frac{1}{(n^2 - \frac{1}{2})} \left(\mu_\lambda(A, \lambda, v)^2 - \frac{|\lambda|^2}{2} \right),$$

as claimed in the lemma. The second statement in the lemma is proved in a very similar fashion. This time we have

$$\begin{aligned} \mathbb{E}_{\dot{A} \in T_A \mathbb{S}} (|\dot{v}|^2) &= \mathbb{E}_{\dot{A} \in T_A \mathbb{S}} (\|A_{\lambda, v}^{-1} P_{v^\perp} \dot{A}v\|^2) = \\ &= \mathbb{E}_{\dot{B} \in \mathcal{N}_{A^\perp}} \mathbb{E}_{t \sim \mathcal{N}(0, \frac{1}{2})} \left(\left\| A_{\lambda, v}^{-1} P_{v^\perp} (t\sqrt{-1}A + \dot{B}) v \right\|^2 \right) = \\ &= \mathbb{E}_{\dot{B} \in \mathcal{N}_{A^\perp}} \left(\left\| A_{\lambda, v}^{-1} P_{v^\perp} \dot{B}v \right\|^2 \right) = \mathbb{E}_{\dot{w} \in \mathcal{N}_{v^\perp}} (\|A_{\lambda, v}^{-1} \dot{w}\|^2) = \|A_{\lambda, v}^{-1}\|_F^2, \end{aligned}$$

the previous to last equality from claim II and the last coming from the fact that for any matrix $B \in \mathbb{C}^{n \times n}$, $\mathbb{E}_{x \sim \mathcal{N}_{\mathbb{C}^n}} \|Bx\|^2 = \|B\|_F^2$ (note the use of Frobenius instead of operator norm in the last equality: that is a crucial point). The last statement of the lemma then follows from claim I. \square

10 Proof of Theorem 2.33

10.1 Proof of (1) and (2) in Theorem 2.33

Note that (2) is trivial. We thus prove (1). The procedure we suggest to choose $\omega \in \Omega_n$ at random is the following (note that each step requires $\mathcal{O}(n^3)$ arithmetic operations or random choices):

1. Choose $B \sim \mathcal{N}_{\mathbb{C}^{(n-1) \times (n-1)}}$ and let U be the Q factor in the QR decomposition of B , then multiply Q by the diagonal matrix with entries $r_{ii}/|r_{ii}|$ where the r_{ii} are the diagonal elements of the R factor. This produces a unitary matrix U uniformly distributed in \mathcal{U}_{n-1} , see for example [39].

2. Choose $\lambda \sim \mathcal{N}_{\mathbb{C}}$ and $M \sim \mathcal{N}_{\mathbb{C}^{(n-1) \times n}}$. Let $H \in \mathcal{U}_n$ be any unitary matrix such that its last column is in $\ker(M)$ (it is trivial to produce such an H by computing the QR decomposition of the matrix whose columns are $\ker(M)$ and the columns of M). Compute Q as the product of the first $n \times (n-1)$ submatrix of H times U . This produces an element with the uniform distribution in the set of $Q \in \mathcal{S}_{n-1}(\mathbb{C}^n)$ such that $(M, Q) \in \mathcal{A}_n$.
3. If $2\Re(\bar{\lambda}\text{tr}(MQ)) > 1 - |\lambda|^2(n-1)$ then discard λ, M, Q and repeat (1) and (2).
4. Choose $w \sim \mathcal{N}_{\mathbb{C}^{n-1}}$.

The only subtle point is that steps (1) and (2) might have to be repeated an arbitrary number of times. The expected number of times that steps (1) and (2) will be repeated is related to C_n defined in (23) by

$$\begin{aligned} \sum_{k=1}^{\infty} \text{Prob}(\text{step } k \text{ is reached}) &= \sum_{k=1}^{\infty} \text{Prob} (2\Re(\bar{\lambda}\text{tr}(MQ)) > 1 - |\lambda|^2(n-1))^{k-1} \\ &= \sum_{k=1}^{\infty} (1 - C_n^{-1})^{k-1} = \frac{1}{(1 - (1 - C_n^{-1}))} = C_n. \end{aligned}$$

10.2 Proof of (3) in Theorem 2.33

We are now prepared for proving (24). Let $\mathbb{1}$ be the characteristic function of the set

$$\{(\lambda, B) : 2\Re(\bar{\lambda}\text{tr}(B)) \leq 1 - |\lambda|^2(n-1)\} \subseteq \mathbb{C} \times \mathbb{C}^{(n-1) \times (n-1)}.$$

From the definition and Fubini's theorem, for any measurable nonnegative function ϕ defined on \mathcal{V} , the expected value $\mathbb{E}_{\omega \sim \Omega_n} (\phi(\psi_n(\omega)))$ equals:

$$\begin{aligned} C_n \mathbb{E}_M \mathbb{E}_{Q:(M,Q) \in \mathcal{A}_n} \left(\mathbb{E}_{\lambda, w} \left(\phi \left(\begin{pmatrix} \lambda & w^* \\ 0 & MQ + \lambda I_{n-1} \end{pmatrix}, \lambda, e_1 \right) \mathbb{1}(\lambda, MQ) \right) \right) = \\ C_n \mathbb{E}_M \mathbb{E}_{Q:(M,Q) \in \mathcal{A}_n} (\alpha(MQ)), \end{aligned}$$

where $\lambda \sim \mathcal{N}_{\mathbb{C}}$, $M \sim \mathcal{N}_{\mathbb{C}^{(n-1) \times n}}$, $w \sim \mathcal{N}_{\mathbb{C}^{n-1}}$ and $\alpha: \mathbb{C}^{(n-1) \times (n-1)} \rightarrow [0, \infty]$ is defined by

$$\alpha(B) = \mathbb{E}_{\lambda, w} \left(\phi \left(\begin{pmatrix} \lambda & w^* \\ 0 & B + \lambda I_{n-1} \end{pmatrix}, \lambda, e_1 \right) \mathbb{1}(\lambda, B) \right).$$

We are then under the hypotheses of Corollary 6.7. Using this result we obtain

$$\mathbb{E}_{w \sim \Omega_n} (\phi(\psi_n(w))) = \frac{C_n}{\Gamma(n)} \mathbb{E}_{B \sim \mathcal{N}_{\mathbb{C}^{(n-1) \times (n-1)}}} (\alpha(B) |\det(B)|^2).$$

With the change of variables $B + \lambda I_{n-1} = D$, which implies

$$\|B\|_F^2 = \|D\|_F^2 + (n-1)|\lambda|^2 - 2\Re(\bar{\lambda}\text{tr}(D)),$$

this last expression equals

$$\frac{C_n}{\Gamma(n)} \mathbb{E}_{\lambda, w, D} \left(\phi \left(\begin{pmatrix} \lambda & w^* \\ 0 & D \end{pmatrix}, \lambda, e_1 \right) \mid \det(D - \lambda I_{n-1})^2 e^{-|\lambda|^2(n-1) + 2\Re(\bar{\lambda}\text{tr}(D))} \mathbb{1}(\lambda, D - \lambda I_{n-1}) \right),$$

where $D \sim \mathcal{N}_{\mathbb{C}^{(n-1) \times (n-1)}}$. Now, note that

$$\mathbb{1}(\lambda, D - \lambda I_{n-1}) \neq 0 \iff e^{-|\lambda|^2(n-1) + 2\Re(\bar{\lambda}\text{tr}(D))} \leq e.$$

We have thus proved

$$\begin{aligned} \mathbb{E}_{w \sim \Omega_n} (\phi(\psi_n(w))) &\leq \frac{e C_n}{\Gamma(n)} \mathbb{E}_{\lambda, w, D} \left(\phi \left(\begin{pmatrix} \lambda & w^* \\ 0 & D \end{pmatrix}, \lambda, e_1 \right) \mid \det(D - \lambda I_{n-1})^2 \right) \\ &\stackrel{(38)}{=} e n C_n \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}} \left(\frac{1}{n} \sum_{\lambda, v: Av = \lambda v} \phi(A, \lambda, v) \right). \end{aligned}$$

This proves claim (3) in Theorem 2.33. \square

We prove the following (non-sharp) bound for the value of C_n .

Lemma 10.1 *With the notations above,*

$$C_n \leq 4n.$$

PROOF. Note that if $0 < |\lambda| \leq (n-1)^{-1/2}$, then for any nonzero $M \in \mathbb{C}^{(n-1) \times n}$ we have

$$\mathbb{P}_Q \left(2\Re(\bar{\lambda}\text{tr}(MQ)) \leq 1 - |\lambda|^2(n-1) \right) \geq \mathbb{P}_Q \left(2\Re(\bar{\lambda}\text{tr}(MQ)) \leq 0 \right) = \frac{1}{2},$$

the last equality coming from the linearity of the trace. We thus have

$$\mathbb{P}_{\lambda, M, Q} \left(2\Re(\bar{\lambda}\text{tr}(MQ)) \leq 1 - |\lambda|^2(n-1) \right) \geq \frac{1}{\pi} \int_{|\lambda| < (n-1)^{-1/2}} \frac{e^{-|\lambda|^2}}{2} d\lambda = \frac{1 - e^{-\frac{1}{n-1}}}{2}.$$

We thus have

$$C_n \leq \frac{2}{1 - e^{-\frac{1}{n-1}}} \leq 4(n-1) \leq 4n,$$

as claimed. \square

11 Proof of Theorem 1.1

Consider the following algorithm.

Algorithm 7 Relative_Error

Input: $\varepsilon \in (0, 1/2)$, $A \in \mathbb{C}^{n \times n}$, $(\zeta, w) \in \mathbb{C} \times \mathbb{C}^n$

Preconditions: (ζ, w) is a certified approximate eigenpair of A with associated eigenpair (λ, v) , $\|w\| = 1$.

```

 $k := 0$ 
 $(\zeta', w') := (\zeta, w)$ 
repeat
   $(\zeta', w') := N_A(\zeta', w')$  (one Newton iteration)
   $k := k + 1$ 
until  $k \geq \log_2 \log_2 \left( \frac{4\|A\|_F}{\varepsilon|\zeta'|} \right)$ 
return  $(\zeta', w')$ 

```

Output: $(\zeta', w') \in \mathbb{C} \times \mathbb{C}^n$

Postconditions: The algorithm halts if $\lambda \neq 0$. In this case, (ζ', w') is an approximate eigenpair of A with associated eigenpair (λ, v) , and $d_S(w', v) \leq \varepsilon$, and moreover $|\zeta' - \lambda| \leq \varepsilon|\lambda|$ and $\text{dist}_A((\zeta, w), (\lambda, v)) \leq c^*/\mu(A, \lambda, v)(1/2)^{2^{2^k-1}}$

By hypothesis,

$$\text{dist}_A((\zeta, w), (\lambda, v)) \leq \frac{c_*}{\mu(A, \lambda, v)} < 1. \quad (58)$$

Hence, $|\zeta - \lambda|/\|A\|_F \leq \text{dist}_A((\zeta, w), (\lambda, v)) < 1$ and the same bound holds with ζ replaced by ζ' at all the iterations of the algorithm (by Definition 2.11). Using that $|\lambda| \leq \|A\|_F$, we deduce that $|\zeta'| \leq 2\|A\|_F$. Hence, at the end of the **repeat** loop the value k satisfies

$$2^{2^k} \geq \frac{4\|A\|_F}{\varepsilon|\zeta'|} \geq \frac{2}{\varepsilon}. \quad (59)$$

This inequality implies that after k iterations of the loop we have, from the definition of approximate eigenpair and the bound (58), that

$$\frac{|\zeta' - \lambda|}{\|A\|_F} \leq \text{dist}_A((\zeta', w'), (\lambda, v)) \leq \frac{c_*}{\mu(A, \lambda, v)} \left(\frac{1}{2} \right)^{2^k-1} \leq \left(\frac{1}{2} \right)^{2^k-1} = \frac{2}{2^{2^k}} \leq \varepsilon. \quad (60)$$

In particular, $d_{\mathbb{S}}(w', v) \leq \varepsilon$ as we wanted. On the other hand, the first inequality in (59) implies

$$2^{2^k-1}|\zeta'| - \|A\|_F \geq \frac{2\|A\|_F}{\varepsilon} - \|A\|_F \geq \frac{\|A\|_F}{\varepsilon}$$

the last since $\varepsilon < 1$. We now use this inequality together with (60) to obtain

$$\frac{|\zeta' - \lambda|}{|\lambda|} \leq \frac{\|A\|_F}{|\lambda|2^{2^k-1}} \leq \frac{\|A\|_F}{\left(|\zeta'| - \frac{\|A\|_F}{2^{2^k-1}}\right)2^{2^k-1}} = \frac{\|A\|_F}{2^{2^k-1}|\zeta'| - \|A\|_F} \leq \varepsilon,$$

i.e., $|\zeta' - \lambda| \leq \varepsilon|\lambda|$.

It remains to show that `Relative_Error` halts provided $\lambda \neq 0$ and to estimate its average running time when A is drawn from $\mathcal{N}_{\mathbb{C}^{n \times n}}$. For this, we note that as soon as

$$k \geq \log_2 \log_2 \left(\frac{8\|A\|_F}{\varepsilon|\lambda|} \right)$$

we shall have (using (60))

$$|\zeta'| \geq |\lambda| - \frac{\|A\|_F}{2^{2^k-1}} \geq |\lambda| - \frac{\varepsilon|\lambda|}{4} = |\lambda|(1 - \varepsilon/4).$$

Therefore, we will also have

$$\log_2 \log_2 \left(\frac{4\|A\|_F}{\varepsilon|\zeta'|} \right) \leq \log_2 \log_2 \left(\frac{4\|A\|_F}{\varepsilon|\lambda|(1 - \varepsilon/4)} \right) \leq \log_2 \log_2 \left(\frac{8\|A\|_F}{\varepsilon|\lambda|} \right) \leq k.$$

Hence, the stopping condition will hold after at most

$$\log_2 \log_2 \left(\frac{8\|A\|_F}{\varepsilon|\lambda|} \right) \leq \log_2 \log_2 \left(\frac{8\|A\|_F\|A^{-1}\|}{\varepsilon} \right)$$

iterations (we have used that $|\lambda|^{-1} \leq \|A^{-1}\|$ for λ is an eigenvalue of A).

We finally estimate the average cost of `Relative_Error`. Since each iteration of the `repeat` loop requires $\mathcal{O}(n^3)$ operations, this cost is at most $\mathcal{O}(n^3)$ times

$$\mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}} \left(\log_2 \log_2 \left(\frac{8\|A\|_F\|A^{-1}\|}{\varepsilon} \right) \right) \leq \log_2 \log_2 \left(\frac{8}{\varepsilon} \mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}} (\|A\|_F\|A^{-1}\|) \right),$$

where we have used Jensen's inequality. Bounds for the expected value of $\|A^{-1}\|$ when $A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}$ are known, see for example [19, Prop. 4.22] which, together with the Cauchy–Schwartz inequality $\mathbb{E}(fg) \leq (\mathbb{E}(f^2)\mathbb{E}(g^2))^{1/2}$, implies

$$\mathbb{E}_{A \sim \mathcal{N}_{\mathbb{C}^{n \times n}}} (\|A\|_F\|A^{-1}\|) \leq \sqrt{n^2 \frac{e(n+1)}{2}} \leq 2n^{3/2}.$$

The statement follows. \square

Acknowledgments. We want to thank Jim Demmel for helpful comments and Froilán Dopico for suggesting us to work out the case of relative error. P. Bürgisser, F. Cucker and M. Shub want to acknowledge the Simons institute where they spent some time while this work was being carried out.

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