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Asymptotic behaviour of a family of gradient algorithms in $\mathbb{R}^d$ and Hilbert spaces

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Abstract. The asymptotic behaviour of a family of gradient algorithms (including the methods of steepest descent and minimum residues) for the optimisation of bounded quadratic operators in $\mathbb{R}^d$ and Hilbert spaces is analyzed. The results obtained generalize those of Akaike (1959) in several directions. First, all algorithms in the family are shown to have the same asymptotic behaviour (convergence to a two-point attractor), which implies in particular that they have similar asymptotic convergence rates. Second, the analysis also covers the Hilbert space case. A detailed analysis of the stability property of the attractor is provided.

1. Introduction

The paper generalizes the results presented in [16] to other optimisation algorithms of the gradient type. We introduce a class of algorithms, called $P$-gradient algorithms, that differ by the choice of the length of the step made in the gradient direction. The class includes in particular the usual steepest-descent algorithm and the method of minimal residues of Krasnosel’kii and Krein [9,10]. We show that for a quadratic function, the worst asymptotic rate of convergence is the same for the whole class of algorithms considered. It is also true that, expressed in the right framework, all the algorithms in the class behave in a very similar fashion. This analysis complements that presented in [1], [13,14] and Chapter 7 of [15] which concerns steepest descent. Moreover, the analysis in [16] directly applies to all algorithms in the class considered, revealing the asymptotic behaviour for bounded quadratic operators not only in $\mathbb{R}^d$ but also in Hilbert spaces. The worst case behaviour exhibited is fundamental “bottom-line” in the

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1 Not all algorithms using the gradient direction belong to that class, which in particular does not include the spectral-gradient algorithm, see [3], proposed by Barzilai and Borwein in [2]. This method, which has been found in particular examples to allow significant improvement over standard steepest descent, see [18], thus requires a separate treatment. The same is true for steepest descent with relaxation or the combination of steepest descent and Barzilai-Borwein methods, as considered in [19].
study of optimisation whose understanding is critical for building more complex and faster algorithms.

The basic idea is renormalisation, as used throughout [15]. The main result in the finite dimension case is that for any algorithm in the class, in the renormalised space one observes convergence to a two-point attractor which lies in the space spanned by the eigenvectors corresponding to the smallest and largest eigenvalues of the matrix \( A \) of the quadratic operator. The proof for bounded quadratic operators in Hilbert space stems from the proof for \( \mathbb{R}^d \) but is considerably more technical. In both cases, as in [1], the method consists of converting the problem to one containing a special type of operator on measures on the spectrum of the operator. The additional technicalities arise from the fact that in the Hilbert space case the measure, which is associated with the spectral measure of the operator, may be continuous. Another important result concerns bounds on convergence rates, named after Kantorovich, see [7]. For all algorithms in the family considered, the actual asymptotic rate of convergence, although satisfying Kantorovich bounds, depends on the starting point and is difficult to predict. This complex behaviour has consequences for the stability of the attractor, which are discussed following the main results.

The family of gradient algorithms we consider, called \( P \)-gradient algorithms, is introduced in Section 2. Renormalisation is presented there, which, together with the monotonic sequences of Section 2.4, forms the core of the analysis to be conducted. The main results are presented in Section 3 first for the case \( \mathcal{H} = \mathbb{R}^d \), then for the Hilbert space case. They rely on the convergence property of successive transformations of a probability measure, which is presented in Section 4. Again, the two cases \( \mathcal{H} = \mathbb{R}^d \) and \( \mathcal{H} \) a Hilbert space are distinguished, the exposition being much simpler in the former case. The stability of attractors is discussed in Section 5 only in the more general case of a Hilbert space, the case \( \mathcal{H} = \mathbb{R}^d \) not allowing for a significant simplification of the presentation. Finally, Section 6 shows the asymptotic equivalence between several rates of convergence of gradient algorithms. All proofs and some important lemmas are collected in an appendix.

2. A family of gradient algorithms

2.1. \( P \)-gradient algorithms

Let \( A \) be a real bounded self-adjoint (symmetric) operator in a real Hilbert space \( \mathcal{H} \) with inner product \((x, y)\) and norm given by \( \|x\| = (x, x)^{1/2} \). Assume that \( A \) is positive, bounded below, and denote its spectral boundaries by \( m \) and \( M \):

\[
m = \inf_{\|x\|=1} (Ax, x), \quad M = \sup_{\|x\|=1} (Ax, x),
\]

with \( 0 < m < M < \infty \). The function to be minimized corresponds to the quadratic form

\[
f(x) = \frac{1}{2} (Ax, x) - (x, y).
\]
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It is minimum at $x^* = A^{-1}y$, its directional derivative at $x$ in the direction $u$ is

$$\nabla_u f(x) = (Ax - y, u).$$

The direction of steepest descent at $x$ is $-g$, with $g = g(x)$ the gradient at $x$, namely $g = Ax - y$. The minimum of $f$ in this direction is obtained for the optimum step-length

$$\gamma = \frac{(g, g)}{(Ag, g)},$$

which corresponds to the usual steepest-descent algorithm. One iteration of the steepest descent algorithm is thus

$$x_{k+1} = x_k - \frac{(g_k, g_k)}{(Ag_k, g_k)}g_k,$$

(2)

with $g_k = Ax_k - y$ and $x_0$ some initial element in $\mathcal{H}$. We define more generally the following class of algorithms.

**Definition 1.** Let $P(\cdot)$ be a real function defined on $[m, M]$, infinitely differentiable, with Laurent series

$$P(z) = \sum_{-\infty}^{\infty} c_k z^k, \quad c_k \in \mathbb{R} \text{ for all } k,$$

such that $0 < \sum_{-\infty}^{\infty} c_k a^k < \infty$ for $a \in [m, M]$. The $k$-th iteration of a $P$-gradient algorithm is defined by

$$x_{k+1} = x_k - \gamma_k g_k$$

(3)

where the step-length $\gamma_k$ minimizes $(P(A)g_{k+1}, g_{k+1})$ with respect to $\gamma$, with $g_{k+1} = g(x_{k+1}) = g(x_k - \gamma g_k)$.

Direct calculation gives

$$\gamma_k = \frac{(P(A)Ag_k, g_k)}{(P(A)A^2g_k, g_k)}$$

(4)

Note that $AP(A) = P(A)A$ and that the denominator and numerator of $\gamma_k$ are linear in $P(A)$. Also, $\gamma_k$ is scale-invariant in $P(A)$ and $\gamma_k \in [1/M, 1/m]$.

Taking $P(A) = A^{-1}$ gives the steepest-descent algorithm. Choosing $P(A) = I$, the identity operator, is equivalent to choosing the step-length that minimizes the norm of the gradient $g_{k+1}$ at the next point. We then obtain the method of minimal residues introduced in [10] for the solution of linear equations. For any fixed $\alpha \in (0, 1)$, choosing $\gamma_k$ that minimizes $\alpha f(x_k - \gamma g_k) + (1 - \alpha)(g(x_k - \gamma g_k), g(x_k - \gamma g_k))$ with respect to $\gamma$ also gives an algorithm in the family. More generally, we show below how to construct $P$-gradient algorithms, with $P(\cdot)$ a polynomial in $A$, using evaluations of $f(\cdot)$ and $g(\cdot)$ only.
2.2. Practical construction when $P$ is a polynomial

We consider the case where $P(A) = A^q$ for some integer $q \geq -1$. (As mentioned, the cases $q = -1$ and $q = 0$ respectively correspond to the methods of steepest-descent and minimal residues.) The extension to $P(\cdot)$ polynomial in $A$ is straightforward (including also linear combinations with $A^{-1}$), using (3).

The minimisation of $(P(A)g_{k+1}, g_{k+1})$, or the calculation of $\gamma_k$ in (3), requires the calculations of terms of the form $(A^n g, g)$, with $n = q$ or $n = q + 1, q + 2$. As shown below, they are easily obtained from evaluations of $g(\cdot)$ at different points. Notice that this construction implies that one iteration of the algorithm will require several evaluations of $g(\cdot)$. The construction proposed below is not necessarily the most economical one, and evaluations of $f(\cdot)$ and $g(\cdot)$ at different points could be combined to provide more efficient evaluations of terms $(A^n g, g)$.

Our objective here is simply to show that the family of algorithms considered in the paper is not of purely theoretical interest, and that other algorithms than the steepest-descent and minimal residues could also be considered in practice.

Let $(A^n g, g)$ be the term to be evaluated, $n \geq 1$, with $g = g(x)$ the gradient at the current point $x$. Define $x(0) = x$ and

$$x(i+1) = x(i) - \beta g(x(i)),$$

with $\beta$ a fixed positive number (for instance, $\beta$ can be taken equal to the value of $\gamma$ at previous iteration of the algorithm). We obtain

$$g^{(i)} = g(x^{(i)}) = (I - \beta A)^i g.$$

Define $P_i = (g, g^{(i)}) = (g, (I - \beta A)^i g)$. In matrix notation, $P_n = Q_n G_n$, where

$$P_n = (P_0, P_1, \ldots, P_n)^\top, \quad G_n = ((g, g), (Ag, g), \ldots, (A^n g, g))^\top$$

and the entries of the $(n + 1) \times (n + 1)$ matrix $Q_n$ are the binomial coefficients,

$$Q_n = \begin{pmatrix}
1 \\
1 & -\beta \\
1 & -2\beta & \beta^2 \\
1 & -3\beta & 3\beta^2 & -\beta^3 \\
\vdots & \vdots & \vdots & \vdots & \vdots
\end{pmatrix}.$$  

The value of $(A^n g, g)$ is then directly obtained from $G_n = Q_n^{-1} P_n$. The entries of $P_n$, defined by $P_i = (g, g^{(i)})$, are also obtained more economically from

$$P_{2j} = (g^{(j)}, g^{(j)})$$

and

$$P_{2j+1} = (g^{(j+1)}, g^{(j)}).$$

Therefore, the evaluation of $\gamma_k = (P(A)Ag_k, g_k)/(P(A)A^2g_k, g_k)$, with $P(\cdot)$ a polynomial of degree $q$, requires $[q/2] + 2$ gradient evaluations (including the one at $x(0) = x_k$).
2.3. Renormalisation

We can rewrite the iteration (3) as

\[(x_{k+1} - x^*) = (x_k - x^*) - \gamma_k g_k,\]

with \(g_k = g(x_k) = A(x_k - x^*),\) so that

\[g_{k+1} = g_k - \gamma_k A g_k = g_k - \left(\frac{(P(A)Ag_k, g_k)}{(P(A)Ag_k, g_k)}\right) Ag_k.\]

Define the renormalised variable

\[z(x) = \frac{Bg(x)}{(P(A)Ag(x), g(x))^{1/2}},\] (5)

with \(B = [P(A)A]^{1/2}\), the positive square-root of \(P(A)A\), so that \((z(x), z(x)) = 1\). Also define \(z_k = z(x_k)\),

\[\mu^k_j = (A^j z_k, z_k), \ j \in \mathbb{Z},\] (6)

so that \(\mu^k_0 = 1\) for any \(k\) and \(\gamma_k = \mu^k_0/\mu^k_1 = 1/\mu^k_1\). We obtain

\[z_{k+1} = \frac{Bg_{k+1}}{(P(A)Ag_{k+1}, g_{k+1})^{1/2}} = \frac{(I - \gamma_k A)Bg_k}{(I - \gamma_k A)z_k}, \] (7)

that is,

\[z_{k+1} = \frac{(I - A/\mu^k_1)z_k}{(\mu^k_2/\mu^k_1)^2 - 1)^{1/2}}.\]

This gives the updating formula for the moments

\[\mu^k_{j+1} = (A^j z_{k+1}, z_{k+1}) = \frac{\mu^k_j - 2\mu^k_{j+1}/\mu^k_1 + \mu^k_{j+2}/(\mu^k_1)^2}{\mu^k_2/(\mu^k_1)^2 - 1}.\] (8)

In the special case where \(H = \mathbb{R}^d\) we can assume that \(A\) is already diagonalised, with eigenvalues \(0 < \lambda_1 \leq \lambda_2 \leq \cdots \leq \lambda_d\). We can then consider \([z_k]_i^2\), the \(i\)-th component of \(z_k\), as a mass on the eigenvalue \(\lambda_i\), with \(\sum_{i=1}^d [z_k]_i^2 = \mu^k_0 = 1\). Define the discrete probability measure \(\nu_k\) supported on \(\lambda_1, \ldots, \lambda_d\) by \(\nu_k(\lambda_i) = [z_k]_i^2\), so that its \(j\)-th moment is \(\mu^k_j\), \(j \in \mathbb{Z}\). We can then interpret (7) as a transformation \(\nu_k \rightarrow \nu_{k+1}\). The asymptotic behaviour of the sequence \((z_k)\) generated by (7) was studied in [1], see also [5] and Chapter 7 of [13]. The main result is that, assuming \(0 < \lambda_1 < \lambda_2 \leq \cdots \leq \lambda_{d-1} < \lambda_d\), the sequence \((z_k)\) converges to a two-dimensional plane, spanned by the eigenvectors \(e_1, e_d\) associated with \(\lambda_1\) and \(\lambda_d\). The attraction property is stated more precisely in Section 3 also in the Hilbert space case. It is already important to notice that although the results in the references above were obtained for the steepest-descent algorithm, the renormalisation [5], which depends on the chosen \(P(.)\), makes them applicable to any algorithm in the family considered. Also, using the renormalisation just defined we easily obtain (non asymptotic) results on the monotonicity of the algorithm along its trajectory.
2.4. Monotonicity of a rate of convergence

Consider the function \((P(A)g_{k+1}, g_{k+1})\) that \(\gamma_k\) minimizes, and compute the rate of convergence \(r_k\) of the algorithm at iteration \(k\), defined by

\[
r_k = \frac{(P(A)g_{k+1}, g_{k+1})}{(P(A)g_k, g_k)}. \tag{9}
\]

Other rates of convergence will be considered in Section 6 where they will be shown to be asymptotically equivalent to \(r_k\). Direct calculation gives

\[
L_k = \frac{\mu_i^k}{\mu_i^{k-1}},
\]

where the moments \(\mu_i^k\) are defined by (6). Also, from (8), \(L_k\) satisfies

\[
L_{k+1} - L_k = \frac{\mu_i^k}{D_k^2} \det M_k
\]

with

\[
D_k = \mu_2^k - (\mu_1^k)^2
\]

and

\[
M_k = \begin{pmatrix}
\mu_{i-1}^k & \mu_i^k & \mu_1^k \\
\mu_0^k & \mu_2^k & \mu_1^k \\
\mu_1^k & \mu_2^k & \mu_3^k
\end{pmatrix}.
\tag{10}
\]

The moment matrix \(M_k\) is positive semi-definite so that \(\det M_k \geq 0\), and thus \(L_{k+1} \geq L_k\), that is, both \(L_k\) and the rate \(r_k\) are non-decreasing along the trajectory followed by the algorithm. When \(\mathcal{H} = \mathbb{R}^2\) (\(d = 2\)), \(\det M_k = 0\) and \(r_k\) is constant. When \(d > 2\) or \(\mathcal{H}\) is a Hilbert space, the rate is monotonically increasing for a typical \(x_0\), indeed, for almost all \(z_0 = z(x_0)\) with respect to the uniform measure on the unit sphere when \(\mathcal{H} = \mathbb{R}^d\). Notice that if the rate is constant over two iterations (\(\det M_k = 0\)), then the measure \(\nu_k\) is supported on two points only, and the iteration (11) for the masses shows that this situation will continue: the rate will thus remain constant for all subsequent iterations.

Note that \(L_k\) and \(D_k\) are bounded (since \(\nu_k\) has a bounded support), respectively by \(L^*\) and \(D^*\), with \(L^* = (M + m)^2/(4mM)\) and \(D^* = (M - m)^2/4\), see Lemma \(\text{(III)}\) in Appendix A3. Therefore, since \(L_k\) is non-decreasing it converges to some limit, and

\[
\det M_k = \frac{(L_{k+1} - L_k)D_k^2}{\mu_1^k} \leq \frac{(L_{k+1} - L_k)(D^*)^2}{m} \to 0, \quad k \to \infty. \tag{11}
\]

In addition to \(L_k\) and \(r_k\) another quantity also turns out to be non-decreasing along the trajectory. Consider

\[
\frac{(P(A)Ag_{k+1}, g_{k+1})}{(P(A)Ax_{k+1} - x_k), (x_{k+1} - x_k)} = \frac{(P(A)Ag_{k+1}, g_{k+1})}{\gamma_k^2 (P(A)Ag_k, g_k)} = \mu_2^k - (\mu_1^k)^2 = D_k. \tag{12}
\]
Direct calculation using (7) gives

\[ D_{k+1} - D_k = \frac{1}{D_k} \det N_k \]

with

\[ N_k = \begin{pmatrix} 
\mu_0^k & \mu_1^k & \mu_2^k \\
\mu_1^k & \mu_2^k & \mu_3^k \\
\mu_2^k & \mu_3^k & \mu_4^k 
\end{pmatrix}. \]

Again, \( N_k \) is positive semi-definite and \( \det N_k \geq 0 \) so that \( D_k \) is also non-decreasing. It converges to some limit and \( \det N_k \) converges to zero for the same reasons as above.

Substitution of \( P(A) \) for a particular algorithm shows which quantities are monotonic. For the steepest-descent algorithm, \( P(A) = A^{-1}, (A^{-1} g_k, g_k) = 2[f(x_k) - f(x^*)] \), and thus the ratios \( r_k = [f(x_{k+1}) - f(x^*)]/[f(x_k) - f(x^*)] \) and \( D_k = (g_{k+1}, g_{k+1})/((x_{k+1} - x_k), (x_{k+1} - x_k)) \) are monotonically non-decreasing. For the method of minimal residues, \( P(A) = I \), and the ratios \( r_k = (g_{k+1}, g_{k+1})/ (g_k, g_k) \) and \( D_k = (Ag_{k+1}, g_{k+1})/ (A(x_{k+1} - x_k), (x_{k+1} - x_k)) \) are monotonically non-decreasing.

The monotonicity and boundedness of \( L_k \) and \( D_k \) makes them suitable for studying the asymptotic behaviour of the algorithm. This is developed in the next section.

### 3. Asymptotic behaviour of gradient algorithms

Consider the case \( \mathcal{H} = \mathbb{R}^d \), and assume that the minimal and maximal eigenvalues of \( A \), \( \lambda_1 = m \), \( \lambda_d = M \), are simple. The attraction property can be stated as follows. Choose \( z_0 = z(x_0) \), the renormalised variable defined by (5) at the initial point \( x_0 \), such that \( (z_0, e_1) > 0 \), \( (z_0, e_d) > 0 \), with \( e_1 \) and \( e_d \) the eigenvectors associated with \( \lambda_1 \) and \( \lambda_d \) respectively. Then

\[ z_{2k} \rightarrow \sqrt{p} e_1 + \sqrt{1 - p} e_d, \quad z_{2k+1} \rightarrow \sqrt{1 - p} e_1 - \sqrt{p} e_d \text{ when } k \rightarrow \infty, \]

where \( p \) is some number in \((0, 1)\), see Section 4 concerning the range of possible values for \( p \). This property, stated in a more general framework in Theorem 1 below, has important consequences for the asymptotic rate of convergence of the algorithm, see Section 6. The proof of the attraction property relies on the convergence of successive transformations of the probability measures \( \nu_k \) defined by \( [z_k]^2 \). The approaches used in [15] to study this convergence do not apply when \( \mathcal{H} \) is infinite dimensional, and we shall present a more general proof in Section 4. It differs somewhat from the one in Chapter 7 of [15], in particular in the choice of the monotonic sequence, \( (L_k) \) instead of \( (D_k) \).

The attraction theorem in \( \mathbb{R}^d \) can be stated as follows. We can assume that \( A \) is diagonalised, and the probability measure \( \nu_k \) is then discrete and puts mass \( [z_k]^2 \) at the eigenvalue \( \lambda_i \). Notice that the updating rule (7) is identical for \( [z_k]^2 \) and \( [z_k] \) associated with \( \lambda_i = \lambda_j \), and the corresponding masses can thus be summed. We can therefore assume that all eigenvalues are different when studying the evolution of \( \nu_k \), see Theorem 3.
**Theorem 1.** Let \( A \) be a \( d \times d \) symmetric matrix, positive definite, with minimum and maximum eigenvalues \( m \) and \( M \) such that \( 0 < m < M < \infty \) and apply a \( P \)-gradient algorithm, see Definition 1 for the minimisation of \( f(x) \) given by (7), initialized at \( x_0 \), with \( z_0 = z(x_0) \), see (4). Assume that

\[
E_1 z_0 \neq 0 \text{ and } E_d z_0 \neq 0,
\]

where \( E_1 \) and \( E_d \) are the orthogonal projectors on the eigenspaces respectively associated with \( \lambda_1 = m \) and \( \lambda_d = M \). Then the asymptotic behaviour of the renormalised gradient \( z_k = z(x_k) \) is such that

\[
z_{2k} = \sqrt{p} u_{2k} + \sqrt{1-p} v_{2k}, \quad z_{2k+1} = \sqrt{1-p} v_{2k+1} - \sqrt{p} v_{2k+1},
\]

with \( \|u_n\| = \|v_n\| = 1 \) \( \forall n \), \( \|Av_n - Mu_n\| \to 0 \), \( \|Av_n - Mv_n\| \to 0 \) as \( n \to \infty \), and \( p \), some number in \((0, 1)\), depending on \( z_0 \).

The proof is omitted since we prove later a more general property valid for \( \mathcal{H} \) a Hilbert space. A more precise result is obtained when the eigenvalues \( \lambda_1 \) and \( \lambda_d \) are simple: the vector \( z_d \) converges to the two-dimensional plane defined by the eigenvectors \( e_1 \) and \( e_d \) associated with \( \lambda_1 \) and \( \lambda_d \).

**Corollary 1** Let \( A \) be a positive-definite symmetric matrix with ordered eigenvalues

\[
0 < m = \lambda_1 < \lambda_2 \leq \cdots \leq \lambda_{d-1} < \lambda_d = M
\]

and let \( e_1, e_d \) be the eigenvectors associated with \( \lambda_1 \) and \( \lambda_d \) respectively. Apply a \( P \)-gradient algorithm, see Definition 1 for the minimisation of \( f(x) \) given by (7), initialized at \( x_0 \) such that \( z_0^\top e_1 \neq 0 \) and \( z_0^\top e_d \neq 0 \), with \( z_0 = z(x_0) \), see (4). Then the algorithm attracts to the plane \( \Pi \) spanned by \( e_1 \) and \( e_d \) in the following sense:

\[
w^\top z_k \to 0, \quad k \to \infty
\]

for any nonzero vector \( w \in \Pi^\perp \). Moreover, the sequence \((z_k)\) converges to a two-point cycle.

This corollary is a straightforward consequence of Theorem 1 when \( \lambda_1 \) and \( \lambda_d \) are simple, with associated eigenvectors \( e_1 \) and \( e_d \), \( u_n \) and \( v_n \) then respectively tend to \( e_1 \) and \( e_d \). The result easily generalizes to the case when (13) is not satisfied. The algorithm then attracts to a two-dimensional plane defined by the eigenvectors \( e_1 \) and \( e_d \) associated with the smallest and largest eigenvalues such that \( z_0^\top e_1 \neq 0 \) and \( z_0^\top e_d \neq 0 \).

We state now the attraction theorem in the more general case where \( \mathcal{H} \) is a Hilbert space. The proof is given in Appendix A1.

**Theorem 2.** Let \( A \) be a bounded real symmetric operator in a Hilbert space \( \mathcal{H} \), positive, with bounds \( m \) and \( M \), such that \( 0 < m < M < \infty \) and apply a \( P \)-gradient algorithm, see Definition 1 for the minimisation of \( f(x) \) given by (7), initialized at \( x_0 \), with \( z_0 = z(x_0) \), see (4). Assume that \( z_0 \) is such that for any \( \epsilon, 0 < \epsilon < (M-m)/2 \),

\[
(E_{m+\epsilon} z_0, z_0) > 0 \text{ and } (E_{M-\epsilon} z_0, z_0) < 1,
\]

(14)
with \((E_\lambda)\) the spectral family of projections associated with \(A\). The asymptotic behaviour of the renormalised gradient \(z_k = z(x_k)\) is such that
\[
z_{2k} = \sqrt{p} u_{2k} + \sqrt{1-p} v_{2k}, \quad z_{2k+1} = \sqrt{1-p} u_{2k+1} - \sqrt{p} v_{2k+1},
\]
with \(\|u_n\| = \|v_n\| = 1 \forall n, \|Au_n - mu_n\| \to 0, \|Av_n - Mv_n\| \to 0\) as \(n \to \infty\), and \(p\), some number in \((0,1)\), depending on \(z_0\).

4. A property of successive transformations of a probability measure

The two properties established in this section form the cornerstones of the proofs of the theorems of previous section. We consider first the case of a discrete measure with finite support, which in terms of convergence of a \(P\)-gradient algorithm corresponds to the case \(\mathcal{H} = \mathbb{R}^d\). The proof is given in Appendix A2.

**Theorem 3.** Let \(\nu_0\) be a discrete probability measure on \(\{\lambda_1, \ldots, \lambda_d\}\) with
\[
0 < m = \lambda_1 < \lambda_2 < \cdots < \lambda_{d-1} < \lambda_d = M < \infty.
\]
Let \([z_k]_i^2\) denote the mass placed at \(\lambda_i\) by \(\nu_k\), that is, \(\nu_k(\lambda_i) = [z_k]_i^2\). Consider the transformation \(T : \nu_k \to \nu_{k+1}\) defined by
\[
[z_{k+1}]_i^2 = \frac{(1 - \lambda_i/\mu_{k}^{i})[z_k]_i^2}{(\mu_{k}^{i}/(\mu_{k}^{i})^2 - 1)^{1/2}}
\]
with the moments \(\mu_{k}^{i}\) defined by (7). Then, when \(k \to \infty\),
\[
([z_{2k}]_1^2) \to p, \quad ([z_{2k+1}]_1^2) \to 1 - p \quad \text{and} \quad ([z_{2k}]_d^2) \to 1 - p, \quad ([z_{2k+1}]_1^2) \to p \tag{17}
\]
for some \(p\) depending on \(\nu_0\), \(0 < p < 1\). Furthermore,
\[
p = \frac{1}{2} \pm \frac{\rho + 1}{\rho - 1} \sqrt{4 - \frac{\rho L}{(\rho + 1)^2}},
\]
with \(\rho = M/m\) and \(L = \lim_{k \to \infty} \mu_{k}^{b} / \mu_{k-1}^{b}\).

Note that the limiting value \(L\) depends on \(\nu_0\), so that the value of \(p\) that characterizes the attractor is difficult to predict. The range of possible values for \(p\) is discussed in Section 5.

We consider now the case of an arbitrary measure on an interval, which raises some additional difficulties compared to previous case. In terms of convergence of a \(P\)-gradient algorithm, it corresponds to the case where \(\mathcal{H}\) is a Hilbert space: for \(E_\lambda\) the spectral family associated with the operator \(A\), we define the measure \(\nu_k\) by \(\nu_k(d\lambda) = d(E_\lambda z_k, z_k), m \leq \lambda \leq M\).
5. Stability of attractors

The range of possible values for $p$ in the attraction Theorem 1 ($\mathcal{H} = \mathbb{R}^d$) is considered in Theorem 3 of [1] (see also Lemma 3.5 of [14]). Let $s(\lambda)$ and $\lambda^*$ be defined by (20). This theorem states that when $\lambda^*$ is not discarded at any iteration, that is, when $\mu^*_k \neq \lambda^*$ for any $k$, then $p \in [1/2 - s(\lambda^*), 1/2 + s(\lambda^*)]$ (note that this assumption cannot be checked). In this section we extend this result in two directions: (i) we will assume that $\mathcal{H}$ is a Hilbert space, (ii) we study the stability of the attractor defined by $p$ in Theorem 2. We shall use the following definition of stability, see [3] p. 444, [11], p. 7.

**Definition 2.** A fixed point $\nu^*$ for a mapping $T(\cdot)$ on a metric space with distance $d(\cdot, \cdot)$ will be called stable if for all $\epsilon > 0$, $\exists \alpha > 0$ such that for any $\nu_0$ for which $d(\nu_0, \nu^*) < \alpha$, $d(T^n(\nu_0), \nu^*) < \epsilon$ for all $n > 0$. A fixed point $\nu^*$ is unstable if it is not stable.

We shall use the distance $d(\nu, \nu')$ given by the Lévy-Prokhorov metric, see [20] p. 349. In our case (measures supported on $[m, M]$), $d(\nu, \nu')$ becomes the Lévy distance between the distribution functions $F, F'$ associated with $\nu, \nu'$, which we denote

$$L(F, F') = \inf \{ \epsilon : F'(x - \epsilon) - \epsilon \leq F(x) \leq F'(x + \epsilon) + \epsilon, \forall x \}.$$ 

In the case where one of the two measures is the discrete measure $\nu^*_p$ concentrated on $m, M$, with $\nu^*_p(m) = p, \nu^*_p(M) = 1 - p$, we get

$$d(\nu, \nu^*_p) = L(F, F^*_p)$$

$$= \inf \{ \epsilon : F(x) \leq p + \epsilon \text{ for } x < M - \epsilon \text{ and } p - \epsilon \leq F(x) \text{ for } m + \epsilon \leq x \},$$

with $F^*_p$ the distribution function associated with $\nu^*_p$. We have then proved the following, see Appendix A4.
Theorem 5. Consider the situation of Theorem 4 with $\nu_0$ any probability measure supported on some closed subset $SS_A$ of $[m, M]$ and
\[
\text{ess inf}(\nu_0) = m, \quad \text{ess sup}(\nu_0) = M.
\]
(i) The measure $\nu_p^*$ is a fixed point for the mapping $T^2$.
(ii) Consider the set $I_u$ defined by
\[
I_u = \left(0, \frac{1}{2} - s(\lambda^*)\right) \cup \left(\frac{1}{2} + s(\lambda^*), 1\right),
\]
where
\[
s(\lambda) = \sqrt{(M - \lambda)^2 + (\lambda - m)^2} \quad \text{and} \quad \lambda^* = \min_{\lambda \in SS_A} s(\lambda). \tag{20}
\]
Any fixed point $\nu_p^*$ with $p$ in $I_u$ corresponds to an unstable fixed point for $T^2$.
(iii) Any point in the interval
\[
I_s = \left(\frac{1}{2} - s(\lambda^*), \frac{1}{2} + s(\lambda^*)\right) \tag{21}
\]
corresponds to a stable $\nu_p^*$ for the mapping $T^2$.

Remark 1. The convergence $d(\nu_k, \nu_p^*) \to 0$ is equivalent to weak convergence $\nu_k \xrightarrow{w} \nu_p^*$ in the usual sense. If $z_k$ is associated with the spectral measure $\nu_k$ and $z_p^*$ with $\nu_p^*$, then, in the Hilbert space this is equivalent to $(z_k - z_p^*, y) \to 0$ for any $y \in \mathcal{H}$, whereas strong convergence would require $\|z_k - z_p^*\| \to 0$. For $\mathbb{R}^d$, the two types of convergence are equivalent, and thus Corollary 1 implies strong convergence. However, for $\mathcal{H}$ a Hilbert space the equivalence is false, and indeed strong convergence generally does not hold. The stability property (iii) is thus a weak statement when $\mathcal{H}$ is a Hilbert space. The $L_2$ metric in $\mathcal{H}$ induces the Hellinger metric on the space of spectral measures, which defines the same topology as the distance in variation, see [20], p. 364. Strong convergence in $\mathcal{H}$ is thus related to distance in variation in the space of spectral measures and is clearly difficult to obtain — except in the special situation where $\nu_0$ has positive mass at $\{m\}$ and $\{M\}$ and presents a spectral gap: $\nu_0([m, m + \epsilon]) = 0$ and $\nu_0([M - \epsilon, M]) = 0$ for some $\epsilon > 0$.

We have $\nu_{k+2}(d\lambda) = H(\nu_k, \lambda)\nu_k(d\lambda)$, with $H(\nu_k, \lambda)$ given by (29) in Appendix A4. One may then notice that when $\nu_0$ is a discrete probability measure, the condition $H(\nu_p^*, \lambda) > 1$ used in the proof of the instability part of the theorem, see Appendix A4, corresponds to a condition on the eigenvalues of the Jacobian of the transformation $T^2$, see [15].

Note that the stability interval $I_s$ always contains the interval
\[
\left(\frac{1}{2} - \frac{1}{2\sqrt{2}}, \frac{1}{2} + \frac{1}{2\sqrt{2}}\right) \approx (0.14645, 0.85355).
\]
Numerical simulations for $\mathcal{H} = \mathbb{R}^3$, with $A$ having eigenvalues $m < \lambda < M$, show that for any initial density of $x_0$ in $\mathbb{R}^d$ associated with a density of $z_0$ reasonably
spread on the unit sphere, the density of the values of $p$ corresponding to stable attractors $\nu_p^*$ can be approximated by

\[
\varphi(p) = C \log[\min\{1, H(\nu_p^*, \lambda)\}] = \begin{cases} 
C \log H(\nu_p^*, \lambda) & \text{if } p \in I_s \\
0 & \text{otherwise,}
\end{cases} \tag{22}
\]

where $C$ is a normalisation constant and $H(\nu_p^*, \lambda)$ is given by (30). Figure 1 shows the empirical density of attractors (full line) together with $\varphi(p)$ (dashed line) in the case $m = 1, \lambda = 4, M = 10$. The support of this density coincides with the stability interval $I_s$ given by (21). When $d > 3$, the density of attractors depends on the initial density of $x_0$.

6. Rates of convergence

We first state a property showing that different definitions of rates of convergence are asymptotically equivalent, see Appendix A5 for the proof.

**Theorem 6.** Let $W$ be a bounded positive self-adjoint operator in $\mathcal{H}$, with bounds $c$ and $C$ such that $0 < c < C < \infty$. Assume that $W$ commutes with $A$ (when $\mathcal{H} = \mathbb{R}^d$, $W$ is a $d \times d$ positive-definite matrix with minimum and maximum eigenvalues respectively $c$ and $C$). Define

\[
R_k(W) = \frac{(W g_{k+1}, g_{k+1})}{(W g_k, g_k)}
\]
if \( \| g_k \| \neq 0 \) and \( R_k(W) = 1 \) otherwise. Apply a \( P \)-gradient algorithm, initialized at \( x_0 \), with \( \gamma_k \) given by (4), for the minimisation of \( f(x) \) given by (1), with minimum value at \( x^* \). Then the limit

\[
R(W,x_0,x^*) = \lim_{n \to \infty} \left[ \prod_{k=0}^{n-1} R_k(W) \right]^{1/n}
\]

exists for all \( x_0, x^* \) in \( \mathcal{H} \) and \( R(W,x_0,x^*) = R(x_0,x^*) \) does not depend on \( W \). In particular,

\[
R(W,x_0,x^*) = \lim_{n \to \infty} \left( \prod_{k=0}^{n-1} r_k \right)^{1/n}
\]

with \( r_k \) defined by (9).

From the results of Section 3, we have

\[
R(W,x_0,x^*) = r(p) = \frac{p(1-p)(\rho - 1)^2}{p + \rho(1-p)||1-p\|p}.
\]

for any \( W \), where \( p \) defines the attractor, see (15), and \( \rho = M/m \) is the condition number of the operator. The function \( r(p) \) is symmetric with respect to \( 1/2 \) and monotonously increasing from 0 to \( 1/2 \), see Figure 2. The worst asymptotic rate is thus obtained at \( p = 1/2 \):

\[
R_{\text{max}} = \left( \frac{\rho - 1}{\rho + 1} \right)^2.
\] (23)

Note that \( \forall k, r_k \leq R_{\text{max}} \) since \( r_k \) is not decreasing, see Section 2.4. For a typical \( x_0 \) (such that the convergence is not finite, that is, such that \( r(p) \neq 0 \)), the stability analysis of Section 5 shows that only values of \( p \) in \( I_s \) given by (21) may correspond to stable attractors. The range of possible values of \( R(p) \) is thus \([R_{\text{min}}, R_{\text{max}}]\), where \( R_{\text{max}} \), given by (23), is obtained for \( p = 1/2 \) and

\[
R_{\text{min}} \leq R_{\text{min}}^* = R(1/2 + 1/[2\sqrt{2}]) = \frac{(\rho - 1)^2}{(\rho + 1)^2 + 4\rho}.
\]

Figure 3 presents the range \([R_{\text{min}}^*, R_{\text{max}}]\) as a function of \( 1/\rho \), the upper curve corresponding to \( R_{\text{max}} \) and the lower to \( R_{\text{min}}^* \). The maximum size of the range is \( 3 - 2\sqrt{2} \approx 0.1716 \), obtained at \( \rho = 1 + 2\sqrt{2} + 2\sqrt{2} + \sqrt{2} \approx 7.5239 \). These results confirm the experimental observation that the rate of convergence of the gradient algorithm is generally close to its worst value \( R_{\text{max}} \), see [14]. The same property is true for any \( P \)-gradient algorithm.

Remark 2. A similar analysis for \( D_k \) defined by (12), which is also not decreasing, shows that \( D_k \to D(p) = p(1-p)(M - m)^2 \) as \( k \to \infty \), with \( D_k \leq D^* = D(1/2) = (M - m)^2/4 \) for all \( k \). Also, for any typical \( x_0 \) such that \( p \in I_s \) given by (21), we have \( D(p) \geq D(1/2 + 1/[2\sqrt{2}]) = (M - m)^2/8 \).
Another quantity of interest is given by
\[
\Delta_N = \log(\frac{R_{\text{max}}}{R_{\text{min}}})/\left[\log(R_{\text{max}}) \log(R_{\text{min}})\right].
\]
Indeed, for \( N \) large enough, \((Wg_N, g_N)/(Wg_0, g_0) \simeq r(p)^N\), the number \( N \) of iterations required for obtaining a ratio \((Wg_N, g_N)/(Wg_0, g_0) = \epsilon (\epsilon \ll 1)\) is approximately \(\log(\epsilon)/\log[r(p)]\) and \(\Delta_N |\log(\epsilon)|\) thus indicates the length of the interval of possible values for \( N \) due to the range of possible values for \( p \). Direct calculation gives \(\Delta_N |\log(R_{\text{max}})| < 1/2\) for any \( p \) and
\[
\Delta_N = \rho/8 - 1/4 + \mathcal{O}(1/\rho), \quad 1/\log(R_{\text{max}}) = -\rho/4 + \mathcal{O}(1/\rho)
\]
for large \( \rho \). Therefore, the number of iterations required by a \( P \)-gradient algorithm to achieve a given precision \( \epsilon \ll 1 \) varies at most by a factor 2 depending on the (typical) starting point \( x_0 \), factors of variation close to 2 being possible only when \( \rho \) is large.

The average value of \( R(W, x_0, x^*) \) for \( z_0 = z(x_0) \) uniformly distributed on the unit sphere is the same for any \( P \)-gradient algorithm, more generally, the distribution of \( R(W, x_0, x^*) \) associated with a particular distribution of \( z_0 \) does not depend on the particular \( P \)-gradient algorithm considered. Moreover, numerical simulations show that the average value of \( R(I, x_0, x^*) \) is the same for the steepest-descent \((P(A) = A^{-1})\) and minimum residues \((P(A) = I)\) algorithms for \( x_0 \) uniformly distributed on the sphere \( \|x_0 - x^*\| = 1 \). The small deviations in average performance between different \( P \)-gradient algorithms can only be related to the fact that a fixed distribution for \( x_0 \) corresponds to different distributions for \( z(x_0) \).

Fig. 2. \( r(p) \) as a function of \( p \), for \( \rho = 2 \) (bottom curve), 4, 8 and 16 (top)
Remark 3. It is known that the introduction of a relaxation coefficient $\gamma$, with $0 < \gamma < 1$, in the steepest-descent algorithm totally changes its behaviour, see, e.g., Chapter 7 of [15]; the algorithm (2) then becomes $x_{k+1} = x_k - \gamma[(g_k, g_k)/(Ag_k, g_k)]g_k$. For $\mathcal{H} = \mathbb{R}^d$ and a fixed $A$, depending on the value of $\gamma$, the renormalized process either converges to periodic orbits (the same for almost all starting points) or exhibits a chaotic behaviour, with the classical period-doubling phenomenon in the case $d = 2$. In higher dimensions, repeated numerical trials show that the process typically no longer converges to the 2-dimensional plane spanned by $(e_1, e_d)$. A detailed analysis for $d = 2$ and experimental results for $d > 2$ also show that relaxation (with $\gamma$ close to 1) considerably improves the rate of convergence. Similar results hold more generally for all $P$-gradient algorithms, with the iteration (3) transformed into $x_{k+1} = x_k - \gamma_k g_k$, with $\gamma$ the (fixed) relaxation coefficient and $\gamma_k$ given by (4). Steepest descent with random relaxation coefficient $\gamma \in (0, 2)$ is considered in [19], avoiding the two point attraction and significantly improving the behavior of ordinary steepest descent.

Appendix

A1. Proof of Theorem 4. The proof relies on Theorem 4 (Theorem 3 when $\mathcal{H} = \mathbb{R}^d$), which concerns successive transformations applied to a probability measure.

Since $A$ is self-adjoint, its spectrum $\text{SS}_A$ is a closed subset of the interval $[m, M]$ of the real line and $m, M \in \text{SS}_A$. Let $E_\lambda$ be the spectral family associated with $A$, and define the spectral measure $\nu_k$ by $\nu_k(d\lambda) = d(E_\lambda z_k, z_k)$, $m \leq \lambda \leq M$. 

![Fig. 3. Range $[R_{\text{min}}^*, R_{\text{max}}]$ of possible values of the asymptotic rate $r(p)$ as a function of $1/\rho$](image)
Since \((2k, z_k) = \int_0^M \nu_k (d\lambda) = 1\), \(\nu_k\) is a probability measure on the Borel sets of \((0, \infty)\), with \(\nu_k([m, M]) = 1\) \(\forall k\). This representation gives

\[
\mu_1 = (Az_k, z_k) = \int \lambda \nu_k (d\lambda), \quad \mu_2 = (A^2 z_k, z_k) = \int \lambda^2 \nu_k (d\lambda)
\]

where integration is over \([m, M]\) unless otherwise specified. Therefore, for any Borel set \(\mathcal{A}\) the transformation (7) gives in terms of \(\nu_k\):

\[
\nu_{k+1} (\mathcal{A}) = \int_{\mathcal{A}} \left[ \lambda - \int \lambda' \nu_k (d\lambda') \right] \nu_k (d\lambda) / \int \lambda^2 \nu_k (d\lambda) = \left[ \int \lambda' \nu_k (d\lambda') \right] / \int \lambda^2 \nu_k (d\lambda).
\]

The conditions (14) on \(\|E\|_{A}\) are equivalent to \(\text{ess inf}(\nu_0) = m\) and \(\text{ess sup}(\nu_0) = M\), see Theorem 4, and the updating rule for \(\nu_k\) can be written as (18). Theorem 4 then implies (19), which can be written as: \(\forall \epsilon > 0, \epsilon \leq \beta = (M - m)/2\),

\[
(E_{m+\epsilon 2k}, z_{2k}) \to p, \quad (E_{M-\epsilon 2k}, z_{2k}) \to p, \quad (E_{m+\epsilon 2k+1}, z_{2k+1}) \to 1 - p, \quad (E_{M-\epsilon 2k+1}, z_{2k+1}) \to 1 - p,
\]

as \(k \to \infty\), where \(p\) depends on \(z_0\), \(0 < p < 1\). Define \(p_{2k} = (E_{m+\beta 2k}, z_{2k})\), \(p_{2k+1} = 1 - (E_{m+\beta 2k+1}, z_{2k+1})\), and the angles \(\phi, \phi_n\) by \(\cos \phi = \sqrt{p}, \sin \phi = \sqrt{1 - p}\), \(\cos \phi_n = \sqrt{p_n}, \sin \phi_n = \sqrt{1 - p_n}\), \(\forall n\). Also define \(s_{2k} = E_{m+\beta 2k} / \cos \phi_{2k}\), \(s_{2k+1} = E_{m+\beta 2k+1} / \sin \phi_{2k+1}\), \(t_{2k} = (z_{2k} - E_{m+\beta 2k}) / \sin \phi_{2k}\), \(t_{2k+1} = -(z_{2k+1} - E_{m+\beta 2k+1}) / \cos \phi_{2k+1}\). This gives \(p_n \to p\) as \(n \to \infty\), \(\|s_n\| = \|t_n\| = 1\) \(\forall n\), and \(z_{2k} = \cos \phi_{2k} s_{2k} + \sin \phi_{2k} t_{2k}\), \(z_{2k+1} = \sin \phi_{2k+1} s_{2k+1} - \cos \phi_{2k+1} t_{2k+1}\). Also,

\[
\|As_n - ms_n\|^2 = \int (\lambda - m)^2 d(E_{\lambda}s_n, s_n),
\]

which, for \(n = 2k\) and any \(\epsilon, 0 < \epsilon < \beta\), gives

\[
\|As_{2k} - ms_{2k}\|^2 = \int_{m}^{m+\epsilon} \frac{(\lambda - m)^2}{p_{2k}} d(E_{\lambda}z_{2k}, z_{2k})
\]

\[
= \int_{m}^{m+\epsilon} \frac{(\lambda - m)^2}{p_{2k}} d(E_{\lambda}z_{2k}, z_{2k}) + \int_{m}^{m+\epsilon} \frac{(\lambda - m)^2}{p_{2k}} d(E_{\lambda}z_{2k}, z_{2k})
\]

\[
\leq \frac{\epsilon^2}{p_{2k}} + \beta^2 \frac{p_{2k} - \int_{m}^{m+\epsilon} d(E_{\lambda}z_{2k}, z_{2k})}{p_{2k}}.
\]

Since \(p_{2k} \to p\) and \(\int_{m}^{m+\epsilon} d(E_{\lambda}z_{2k}, z_{2k}) \to p\) as \(k \to \infty\), \(\|As_{2k} - ms_{2k}\| \to 0\) as \(k \to \infty\). Similarly, \(\|As_{2k+1} - ms_{2k+1}\| \to 0\) as \(k \to \infty\) and \(\|At_n - Mt_n\| \to 0\) as \(n \to \infty\). Consider now

\[
u_n = \cos \phi, \quad s_n = \sin \phi, \quad t_n = 1 \quad \forall n.
\]

Straightforward calculations show that \(\theta_n = \phi_n - \phi\) gives (15) with \(\|u_n\| = \|v_n\| = 1\) \(\forall n\). Also

\[
\|Au_n - mu_n\| \leq \|u_n\|\|As_n - ms_n\| + \sin \theta_n (M - m).
\]
and, since \( \|A s_n - m s_n\| \to 0 \), \( \theta_n \to 0 \) as \( n \to \infty \), \( \|A u_n - m u_n\| \to 0 \) as \( n \to \infty \).
Similarly, \( \|A v_n - M v_n\| \to 0 \) as \( n \to \infty \).

**A2. Proof of Theorem** We first prove that the mass of \( \nu_k \) tends to concentrate on two eigenvalues only. When \( \nu_k \) is non degenerate, \( L_1 > 1 \) from Jensen inequality, and thus, since \( (L_k) \) is non-decreasing, see Section 2.4. \( L_k \geq L_1 > 1 \). Now, from Lagrange identity \( (\sum a_i^2)(\sum b_i^2) = \sum_{i<j}(a_i b_j - a_j b_i)^2 + (\sum a_i b_i)^2 \)

\[
L_k = \left( \sum_{i=1}^{d} \lambda_i [z_k]^2 \right) \left( \sum_{i=1}^{d} [z_k]^2 / \lambda_i \right)
\]

\[
= \sum_{i<j} [z_k]^2 [z_k]^2 \left( \frac{\sqrt{\lambda_i}}{\sqrt{\lambda_j}} - \frac{\sqrt{\lambda_j}}{\sqrt{\lambda_i}} \right)^2 + \left( \sum_{i=1}^{d} [z_k]^2 \right)^2
\]

\[
= \sum_{i<j} [z_k]^2 [z_k]^2 \frac{(\lambda_i - \lambda_j)^2}{\lambda_i \lambda_j} + 1.
\]

Let \( i_k \) and \( j_k \) denote the indices that achieve \( \max_{i<j}[z_k]^2 [z_k]^2 \). We have

\[
L_k \leq [z_k]^2 [z_k]^2 \sum_{i<j} \frac{(\lambda_i - \lambda_j)^2}{\lambda_i \lambda_j} + 1
\]

and thus

\[
[z_k]^2 [z_k]^2 \geq \delta = \frac{L_1 - 1}{\sum_{i<j} (\lambda_i - \lambda_j)^2}.
\]

Moreover, \( [z_k]^2 + [z_k]^2 < 1 \) gives

\[
\delta < [z_k]^2 < 1 - \delta \text{ and } \delta < [z_k]^2 < 1 - \delta.
\]

Consider the matrix \( M_k \) given by (10). Its determinant can be written as

\[
\det M_k = \sum_{i<j<i<k} [z_k]^2 [z_k]^2 \frac{(\lambda_i - \lambda_j)^2 (\lambda_i - \lambda_k)^2 (\lambda_j - \lambda_k)^2}{\lambda_i \lambda_j \lambda_k}
\]

\[
\geq [z_k]^2 [z_k]^2 (\lambda_{i_k} - \lambda_{j_k})^2 \sum_{i \neq i_k, j_k} [z_k]^2 \frac{(\lambda_i - \lambda_{i_k})^2 (\lambda_j - \lambda_{j_k})^2}{\lambda_i \lambda_{i_k} \lambda_{j_k}}
\]

\[
\geq \frac{\delta^6}{N^2} \sum_{i \neq i_k, j_k} [z_k]^2
\]

where

\[
\delta = \min_{i,j} |\lambda_i - \lambda_j|.
\]

Since \( \det M_k \to 0 \) as \( k \to \infty \), see (11), we get \( \sum_{i \neq i_k, j_k} [z_k]^2 \to 0 \) as \( k \to \infty \). The mass thus tends to concentrate on \( \lambda_{i_k}, \lambda_{j_k} \).
Next we prove that $i_k$ and $j_k$ eventually become fixed. From the result above, \( \forall \epsilon > 0, \exists k_\epsilon \) such that \( \sum_{i \neq i_k, j_k} |z_k|^2 < \epsilon \), \( k > k_\epsilon \).

Consider the updating equation (16). We have for any \( i, (\mu^i_k - \lambda_i)^2 \leq (M - m)^2 \). Also, \( D_k = \mu^k_2 - (\mu^k_1)^2 \geq D_0 \), see Section 2.4. This gives for \( i \neq i_k, j_k \) and \( k > k_\epsilon \)

\[
|z_{k+1}|^2 < \epsilon \frac{(M - m)^2}{D_0}.
\]

Taking \( \epsilon_1 = \delta D_0/(M - m)^2 \) we obtain \( |z_{k+1}|^2 < \delta \) for \( i \neq i_k, j_k \) and \( k > k_\epsilon \). Since \( |z_{k+1}|^2 \lambda \) and thus \( (\lambda_{D} - \mu_1)^2 > (\lambda_{j*} - \mu_1)^2 \) for \( k \) large enough. We have

\[
\mu_1^k = \lambda^{*, i}[z_{k,i}]^2 + \lambda^{*, j}[z_{k,j}]^2 + \sum_{i \neq i^*, j^*} \lambda_i |z_k|^2 \leq \lambda^{*, i}[z_{k,i}]^2 + \lambda^{*, j}[z_{k,j}]^2 + \lambda d \sum_{i \neq i^*, j^*} |z_k|^2.
\]

Take \( \epsilon_2 = \min\{\epsilon_1, \delta \delta_{\lambda} \}. \) For \( k > k_\epsilon \) we have

\[
\mu_1^k \leq \lambda^{*, i}[z_{k,i}]^2 + \lambda^{*, j}[z_{k,j}]^2 + \lambda d \epsilon_2 \leq \lambda^{*, i} \delta + \lambda^{*, j} (1 - \delta) + \lambda d \epsilon_2 \leq \lambda^{*, j} - \delta \delta_{\lambda} + \lambda d \epsilon_2 \leq \lambda^{*, j},
\]

and thus \( \lambda d - \mu_1^k > (\lambda^{*, j} - \mu_1^k)^2 \). From (16), this gives for \( k > k_\epsilon \)

\[
\left( \frac{|z_{k+1}|^2}{|z_k|^2} \right) = \frac{(\lambda d - \mu_1^k)^2}{D_k} > \frac{(\lambda^{*, j} - \mu_1^k)^2}{D_k} = \left( \frac{|z_{k+1}|_{j^*}^2}{|z_k|_{j^*}^2} \right)^2.
\]

We arrived at a contradiction since \( |z_{k+1}|^2 \to 0 \) and \( |z_k|^2 \) is bounded from below by \( \delta \). Therefore \( j^* = d \). Similarly, \( i^* = 1 \).

Finally, let \( L \) denote \( \lim_{k \to \infty} L_k \), see Section 2.4. There are only two discrete measures with nonzero masses on \( \lambda_1 \) and \( \lambda_d \) and such that \( \mu_1 \mu_{-1} = L \),

\[
\nu^{(1)} = \left\{ \lambda_1 \begin{array}{ll} \lambda_d & \text{if } p \ 1 - p \end{array} \right\} \quad \text{and} \quad \nu^{(2)} = \left\{ \begin{array}{ll} \lambda_1 & \lambda_d \end{array} \right\}
\]

with

\[
p = \frac{1}{2} - \frac{\rho + 1}{\rho - 1} \sqrt{\frac{1}{4} - \frac{\rho L}{(\rho + 1)^2}}
\]

and \( \rho = M/m \). Direct calculation shows that \( \nu_k = \nu^{(1)} \) gives \( \nu_{k+1} = \nu^{(2)} \), hence the convergence of \( \nu_k \) to the cyclic attractor \( \nu^{(1)} \to \nu^{(2)} \to \nu^{(1)} \to \cdots \). 

\textbf{A3.} The proof of Theorem 4 is more technical than that of Theorem 3 and relies on a series of lemmas stated below.
Lemma 1. Let $\nu$ be any probability distribution on $[m,M]$, $0 < m \leq M < \infty$ with moments $\mu_i = \int \lambda^i \nu(d\lambda)$, $i \in \mathbb{Z}$ ($\mu_0 = 1$). Then,

\begin{align}
\mu_2 - \mu_1^2 &\leq D^* = (M - m)^2/4 \quad \text{(24)} \\
\mu_2 &- \mu_1 \leq L^* = (M + m)^2/(4mM) \quad \text{(25)}
\end{align}

Proof. The proof relies on standard results in experimental design theory, see, e.g., [4, 24]. Consider the two linear regression models $\eta_1(\theta, \lambda) = \theta_0 + \theta_1 \lambda$ and $\eta_2(\theta, \lambda) = \theta_0/\sqrt{\lambda} + \theta_1 \sqrt{\lambda}$, with $\theta_0, \theta_1$ the model parameters and $\lambda$ the design variable, $\lambda \in [m, M]$. A point optimum design (approximate theory) aims at determining a probability measure on $[m, M]$ that maximizes the determinant of the information matrix associated with a particular model, here respectively

$$I_1(\nu) = \begin{pmatrix} \mu_0 & \mu_1 \\ \mu_1 & \mu_2 \end{pmatrix} \quad \text{and} \quad I_2(\nu) = \begin{pmatrix} \mu_0 & \mu_1 \\ \mu_1 & \mu_2 \end{pmatrix}.$$

The function $\log \det I(\nu)$ is concave on the set of probability measures on $[m, M]$, and its maximum is unique. The Kiefer-Wolfowitz General Equivalence Theorem [8] gives a characterization of the measure $\nu^*$ that maximizes $\det I_1(\nu) = \mu_2 - \mu_1^2$ and $\det I_2(\nu) = \mu_1 \mu_2 - 1$. In this case it corresponds to the Kantorovich inequality, see [7] and [12], p. 151. (A full development of this connection is presented in [17].)

Lemma 2. Let $\nu$ be any probability distribution on $[m,M]$, $0 < m \leq M < \infty$. Assume that there exists an interval $\mathcal{I} \subseteq [m, M]$, $|\mathcal{I}| \leq \alpha$ and $\nu(\mathcal{I}) \geq 1 - \epsilon$, $\epsilon \in [0, 1]$. Then, $\text{Var}(\nu) \leq \alpha^2/4 + 2\epsilon M^2$.

Proof. Define $\mu_1 = \int_{[m,M]} \lambda \nu(d\lambda)$, $\mu_2 = \int_{\mathcal{I}} \lambda \nu(d\lambda)$. Then $\mu_1 = \mu_2 + \int_{[m,M] \setminus \mathcal{I}} \lambda \nu(d\lambda)$. Therefore, $\mu_2 \leq \mu_1 \leq \mu_2 + \epsilon M$. We get

$$\text{Var}(\nu) = \int (\lambda - \mu_1)^2 \nu(d\lambda) \leq \int_{\mathcal{I}} (\lambda - \mu_1)^2 \nu(d\lambda) + (M - m)^2 \epsilon \leq \int \mathcal{I} (\lambda - \mu_2)^2 \nu(d\lambda) + (\mu_1 - \mu_2)^2 \nu(\mathcal{I}) + (M - m)^2 \epsilon.$$

Lemma [4] implies $\int \mathcal{I} (\lambda - \mu_2)^2 \nu(d\lambda) \leq \alpha^2/4$ and $(\mu_1 - \mu_2)^2 \leq \epsilon^2 M^2$ gives

$$\text{Var}(\nu) \leq \alpha^2/4 + \epsilon^2 M^2 + M^2 \epsilon \leq \alpha^2/4 + 2\epsilon M^2.$$

Lemma 3. Let $\nu$ be any probability distribution on $[m,M]$, $0 < m \leq M < \infty$. Assume that $\text{Var}(\nu) \leq \epsilon$. Then, there exist an interval $\mathcal{I}$ such that $|\mathcal{I}| \leq \epsilon^{1/4}$ and $\nu(\mathcal{I}) \geq 1 - 4\sqrt{\epsilon}$.

Proof. Take $\mathcal{I} = [\mu_1 - \epsilon^{1/4}/2, \mu_1 + \epsilon^{1/4}/2]$, $\mu_1 = \int \lambda \nu(d\lambda)$, and apply the Chebyshev inequality.
Lemma 4. Let \( \nu \) be any distribution on \([m, M]\), \( 0 < m \leq M < \infty \). Define 
\[
\mu_i = \int \lambda^i \nu(d\lambda)
\]
and
\[
M = \begin{pmatrix}
\mu_{-1} & \mu_0 & \mu_1 \\
\mu_0 & \mu_1 & \mu_2 \\
\mu_1 & \mu_2 & \mu_3
\end{pmatrix}.
\]
Assume that \( L = \mu_{-1} \mu_1 > 1 \) (which, by Jensen’s inequality, holds when \( \nu \) is not degenerate at a single point) and \( \det M < \epsilon \). Then, there exist two intervals \( I_1 \) and \( I_2 \) such that
\[
(i) \quad |I_i| \leq \frac{(M - m)^{3/4}}{m^{3/4}(L - 1)^{3/2}}, \quad i = 1, 2, \quad \nu(I_1) + \nu(I_2) \geq 1 - 4\sqrt{\epsilon} M^{3/2},
\]
\[
(ii) \quad \max_{x \in I_i} |x - \mu_{-1}| > \frac{3(L - 1)m^2}{4(M - m)}, \quad i = 1, 2,
\]
\[
(iii) \quad \text{for } \epsilon < \epsilon_*, \quad \nu(I_i) \geq \frac{m^2(L - 1)}{4M^2}, \quad i = 1, 2,
\]
\[\text{and } \max_{x \in I_1, y \in I_2} |x - y| > m\sqrt{2(L - 1)}.\]

Proof.
(i) Consider the measure \( \nu' \) defined by \( \nu'(A) = (1/\mu_{-1}) \int_A (1/\lambda) \nu(d\lambda) \) for any Borel set \( A \subset [m, M] \), and denote its moments by \( \mu'_i = (1/\mu_{-1}) \int \lambda^{1-i} \nu(d\lambda) = \mu_{i-1}/\mu_{-1}. \) Note that for any Borel set \( A \)
\[
\frac{1}{\mu_{-1}} \nu(A) \leq \nu'(A) \leq \frac{1}{m \mu_{-1}} \nu(A).
\]
We have
\[
M' = \begin{pmatrix}
\mu'_0 & \mu'_1 & \mu'_2 \\
\mu'_1 & \mu'_2 & \mu'_3 \\
\mu'_2 & \mu'_3 & \mu'_4
\end{pmatrix} = M/\mu_{-1}
\]
and thus \( \det M' = \det M/\mu_{-1}^3 \). Also define \( D' = \mu'_1 - (\mu'_1)^2, \) \( a = \sqrt{D'}, \) \( b = (\mu'_1 \mu'_2 - \mu'_1^2)/\sqrt{D'}, \) \( c = a \mu'_2 + b \mu'_1 = (\mu'_2)^2 - \mu'_1^2 \mu'_3)/\sqrt{D'} \) (note that \( a > 0, b > 0 \) and \( c < 0 \)) and \( \eta = F(\zeta) = a\zeta^2 + b\zeta - c, \) with \( \zeta \) having the distribution \( \nu' \). Direct calculation gives \( E'([\eta]) = \int \eta(\zeta) \nu'(d\zeta) = 0 \) and \( \text{Var}'(\eta) = E'(\eta^2) - (E'(\eta))^2 = \det M' \), so that \( \det M < \epsilon \) implies \( \text{Var}'(\eta) < \epsilon' = \epsilon/\mu_{-1}^3. \) From Lemma 3, the interval \( I = [-\epsilon'^{1/2}/2, \epsilon'^{1/2}/2] \) is such that \( \text{Pr}\{\eta \in I\} \geq 1 - 4\sqrt{\epsilon'}. \) Also, from the mean-value theorem, there exist \( \lambda_1 < \lambda_2 \) such that \( \lambda_1 \in [m, M] \) and \( a\lambda_i^2 + b\lambda_i - c = 0, \) \( i = 1, 2. \) Direct calculation gives \( F(\mu'_1) = F(\mu_{-1}) = a(\mu'_1)^2 + b\mu'_1 - c = -(D')^{3/2}, \) and thus
\[
m \leq \lambda_1 < 1/\mu_{-1} < \lambda_2 \leq M.
\]
Take \( \beta = (M - m)(\epsilon')^{1/4}/[2(D')^{3/2}], \) we get
\[
F(\lambda_1 + \beta) < -(\epsilon')^{1/4}/2, F(\lambda_1 - \beta) > (\epsilon')^{1/4}/2, \quad F(\lambda_2 + \beta) > (\epsilon')^{1/4}/2, F(\lambda_2 - \beta) < -(\epsilon')^{1/4}/2, \quad
\]
and \( \nu(I_1) + \nu(I_2) \geq 1 - 4\sqrt{3}M^{3/2} \) when \( I_i = [\lambda_i - \beta, \lambda_i + \beta], i = 1,2 \), with 
\( |I_i| = 2\beta = (M - m)\epsilon^{1/4} \mu_{i-1}^{3/4}/(L - 1)^{3/2} \leq (M - m)\epsilon^{1/4} m^{3/4}/(L - 1)^{3/2} \).

(ii) Define \( y_1 = \mu'_1 - \lambda_1, y_2 = \lambda_2 - \mu'_2 \), so that \( \max_{x \in I_i} |x - \mu_{i-1}| > y_1 \) and \( \max_{x \in I_2} |x - \mu_{-1}| > y_2 \). We have \( F(\lambda) = a(\lambda - \lambda_1)(\lambda - \lambda_2) \) and thus \( y_1 y_2 = -F'(\mu'_1)/a = D' \). Also, \( |y_2 - y_1| < y_1 + y_2 \leq M - m \), so that \( D' > y_i(y_i + M - m), i = 1, 2 \), and thus

\[
y_i > \frac{M - m}{2} \left[ \sqrt{1 + \frac{4D'}{(M - m)^2}} - 1 \right] > \frac{D'}{M - m} \left( 1 - \frac{D'}{(M - m)^2} \right), i = 1, 2.
\]

Lemma \[4\] gives \( D' < (M - m)^2/4 \), so that

\[
y_i > 3D'/[4(M - m)] > 3(L - 1)m^2/[4(M - m)], i = 1, 2.
\]

(iii) Define \( \gamma = \nu'(I_2) \), part (i) implies \( \nu'(I_1) > 1 - 4\sqrt{3} - \gamma \), and from Lemma \[2\]

\[
D' \leq \frac{(M - m)^2 2\sqrt{3}}{4D'^3} + 2(4\sqrt{3} + \gamma)M^2,
\]

which gives

\[
\gamma \geq \frac{D'}{2M^2} - \sqrt{3} \left[ \frac{(M - m)^2}{8(D')^3M^2} + 4 \right],
\]

and thus \( \gamma \geq D'/(4M^2) > (L - 1)m^2/[4M^2] \) for \( \epsilon < \epsilon_1 < [4(D')^8]/[(M - m)^2 + 32D'^3M^2] \), see \(27\).

Define now \( \Delta = \max_{x \in I_1, y \in I_2} |x - y| \). Lemma \[2\] gives \( D' < \Delta^2/4 + 8\sqrt{3}M^2 \), which implies \( \Delta^2 \geq 4D' - 32M^2\sqrt{3} \). Since \( \epsilon < \epsilon_1 \) implies \( \sqrt{3} < D'/16M^2 \), we get \( \Delta^2 > 2(L - 1)m^2 \). □

**Proof of Theorem \[4\]** The proof follows the same lines as that of Theorem \[3\] and is divided into four parts. In (i), we construct sequences of intervals \( I_k = [m_k, m_{k+1}] \) and \( R_k = [M_k - \delta, M_k] \) in which the measure \( \nu_k \) will tend to concentrate. In (ii) we prove that \( R_k \cap R_{k+1} \neq \emptyset \) and in (iii) that the sequence \( M_k \) is non-decreasing. Finally, the limiting behaviour of \( \nu_k \) is derived in (iv).

(i) We have seen in Section \[2.4\] that \( \det M_k \to 0 \) as \( k \to \infty \), with \( M_k \) given by \(10\). Therefore, given \( \epsilon, \exists K_\epsilon \) such that \( \forall k > K_\epsilon, \det M_k < \epsilon \). Define \( L_k = \mu_k^{1/2}\mu_{k-1} \) and note that \( L_k > 1 \) because no \( \nu_k \) is degenerate at a single point. Using Lemma \[4\] for \( \epsilon \) small enough, for any \( k > K_\epsilon \) there exist two intervals \( I_1^k, I_2^k \), with width at most

\[
\delta = \delta(\epsilon) = \frac{(M - m)\epsilon^{1/4}}{m^{3/4}(L_0 - 1)^{3/2}},
\]

and such that \( \nu_k(I_1^k) + \nu_k(I_2^k) \geq 1 - 4\sqrt{3}M^{3/2}, \nu_k(I_1^k) \geq m^2(L_k - 1)/(4M^2), \nu_k(I_2^k) \geq m^2(L_k - 1)/(4M^2) \). Also, \( \max_{x \in I_1^k, y \in I_2^k} |x - y| \geq m\sqrt{2(L_k - 1)} \). Without any loss of generality, assume that \( I_1^k \) is the interval on the left. Define
\[ \mathcal{L}(x) = [x, x + \delta], \mathcal{R}(x) = [x - \delta, x], \]

\[ \lambda_k^L = \text{Arg max}_x \{ \nu_k(\mathcal{L}(x)) \mid \mathcal{L}(x) \cap \mathcal{I}_1^k \neq \emptyset \}, \]

\[ \lambda_k^R = \text{Arg max}_x \{ \nu_k(\mathcal{R}(x)) \mid \mathcal{R}(x) \cap \mathcal{I}_2^k \neq \emptyset \}, \]

and \( m_k = \min \lambda_k^L, M_k = \max \lambda_k^R, \mathcal{L}_k = \mathcal{L}(m_k), \mathcal{R}_k = \mathcal{R}(M_k) \); that is, \( M_k \) is the right endpoint of an interval \( \mathcal{R}_k \), intersecting \( \mathcal{I}_2^k \), with maximum measure, and similarly for \( m_k \) and \( \mathcal{L}_k \). Note that \( \nu_k(\mathcal{L}_k) + \nu_k(\mathcal{R}_k) \geq 1 - 4\sqrt{\epsilon}M^{3/2}, \nu_k(\mathcal{L}_k) \geq m^2(L_k - 1)/(4M^2) \) and \( \nu_k(\mathcal{R}_k) \geq m^2(L_k - 1)/(4M^2) \). The situation is the same for the two sequences of intervals \( (\mathcal{L}_k) \) and \( (\mathcal{R}_k) \), and we concentrate on \( (\mathcal{R}_k) \) in the rest of the proof.

(ii) We show now that \( \mathcal{R}_k \cap \mathcal{R}_{k+1} \neq \emptyset \). Again for \( \epsilon \) small enough \( \mu^k_1 \notin \mathcal{R}_k \) and \( \lambda - \mu^k_1 \geq M_k - \delta - \mu^k_1 \) on \( \mathcal{R}_k \) so that

\[ \nu_{k+1}(\mathcal{R}_k) = \int_{\mathcal{R}_k} \frac{(\lambda - \mu^k_1)^2}{D_k} \nu_k(d\lambda) \geq \frac{\nu_k(\mathcal{R}_k)}{D_k}(M_k - \delta - \mu^k_1)^2 \]

\[ \geq \frac{m^2(L_k - 1)}{4M^2D^*}(M_k - \delta - \mu^k_1)^2 \]

with \( D^* \) the maximum possible value of \( D_k, D^* = (M - m)^2/4 \), see Lemma 1.

By construction, \( \max_{x \in \mathcal{I}_2^k} |x - \mu^k_1| \leq M_k + \delta - \mu^k_1 \), and thus, from Lemma 3

\[ M_k - \mu^k_1 + \delta \geq \frac{3m^2(L_k - 1)}{4(M - m)} \geq \frac{3m^2(L_0 - 1)}{4(M - m)} = C. \] (28)

Choosing \( \epsilon \) such that \( \delta < C/4 \) gives \( M_k - \delta - \mu^k_1 > C/2 \) and thus

\[ \nu_{k+1}(\mathcal{R}_k) \geq \frac{m^2(L_k - 1)}{4M^2D^*} \frac{C^2}{4} \geq \nu^*_R = \frac{9m^6(L_0 - 1)^3}{16M^2(M - m)^4}. \]

Choosing now \( \epsilon \) such that \( 4\sqrt{\epsilon}M^{3/2} < \nu^*_R \) we obtain \( \mathcal{R}_k \cap \mathcal{R}_{k+1} \neq \emptyset \) for any \( k > K_\epsilon \).

(iii) We prove now that the sequence \( (M_k) \) is not decreasing starting at some \( K_\epsilon \) for \( \epsilon \) small enough. Take \( k > K_\epsilon \) and assume that \( M_{k+1} = M_k - \beta, \beta > 0 \). Then note that \( \beta < \delta \) since \( \mathcal{R}_k \cap \mathcal{R}_{k+1} \neq \emptyset \) by (ii) above. Consider the difference

\[ \nu_{k+1}(\mathcal{R}_k) - \nu_{k+1}(\mathcal{R}_{k+1}) = \nu_{k+1}([M_k - \beta, M_k]) - \nu_{k+1}([M_k - \delta - \beta, M_k - \delta]). \]

Assume first that \( \nu_{k+1}([M_k - \delta - \beta, M_k - \delta]) = 0 \), then \( \nu_{k+1}(\mathcal{R}_k) > \nu_{k+1}(\mathcal{R}_{k+1}) \), which is impossible by construction. We can thus consider the following ratio

\[ \frac{\nu_{k+1}([M_k - \beta, M_k])}{\nu_{k+1}([M_k - \delta - \beta, M_k - \delta])} = \frac{\int_{M_k - \beta} M_k (\lambda - \mu^k_1)^2 \nu_k(d\lambda)}{\int_{M_k - \delta - \beta} M_k (\lambda - \mu^k_1)^2 \nu_k(d\lambda)} \]

\[ \geq \frac{(M_k - \beta - \mu^k_1)^2}{(M_k - \delta - \mu^k_1)^2} \frac{\nu_k([M_k - \beta, M_k])}{\nu_k([M_k - \delta - \beta, M_k - \delta])}. \]
Since $M_k - \delta - \mu^k_i \geq C - 2\delta \geq 2\delta$ for $C > 4\delta$, see (28), and $\beta < \delta$, $(M_k - \beta - \mu^k_i)^2 > (M_k - \delta - \mu^k_i)^2$. Also, by construction,

$$0 \leq \nu_k([M_k - \beta, M_k]) - \nu_k([M_k - \delta - \beta, M_k - \delta]) = \nu_k([M_k - \beta, M_k]) - \nu_k([M_k - \delta - \beta, M_k - \delta]).$$

This gives

$$\frac{\nu_{k+1}([M_k - \beta, M_k])}{\nu_{k+1}([M_k - \delta - \beta, M_k - \delta])} > 1.$$ 

Therefore, $\beta > 0$ leads to $\nu_{k+1}(R_k) > \nu_{k+1}(R_{k+1})$, which is impossible. We thus obtain $M_{k+1} \geq M_k$ for $k > K$. 

(iv) Since the sequence $(M_k)$ is non-decreasing and bounded from above (by $M$), it has a limit $M_* \geq M$. The same is true for $m_k$, and $m_k \to m_*$ as $k \to \infty$. We have thus proved that for any $\delta$ small enough and any $k$ larger than some $K_\delta$,

$$\nu_k([M_* - \delta, M_*]) + \nu_k([m_*, m_* + \delta]) \geq 1 - \frac{4M^{3/2}m^{9/2}(L_0 - 1)^3\delta^2}{(M - m)^2}.$$

Assume that $M_* < M$. This would imply $\nu_k([M - \delta, M]) \to 0$ as $k \to \infty$ for $\delta < M - M_*$. On the other hand,

$$\frac{\nu_{k+1}([M - \delta, M])}{\nu_{k+1}([M_* - \delta, M_*])} > \frac{\nu_k([M - \delta, M])}{\nu_k([M_* - \delta, M_*])},$$

which leads to a contradiction since $\nu_k([M - \delta, M])/\nu_k([M_* - \delta, M_*])$ is then increasing and $\nu_k([M_* - \delta, M_*])$ is bounded from below. Therefore, $M_* = M$, and similarly $m_* = m$, with, for $\delta$ small enough and any $k$ larger than some $K_\delta$, $\nu_k([m + \delta, M - \delta]) < 4M^{3/2}m^{9/2}(L_0 - 1)^3\delta^2/(M - m)^2$. Finally, from Helly’s Theorem, see [20], p. 319, from the sequence $(\nu_k)$ we can extract a subsequence $(\nu_{k_i})$ that is weakly convergent, and from the result above the associated limit has necessarily the form $\nu^*_p$, where $\nu^*_p$ is the discrete measure concentrated on the two points $m$, $M$, with $\nu^*_p(m) = p$, $\nu^*_p(M) = 1 - p$. Since $L_k$ converges to some $L$, $\nu^*_p$ is such that the associated value of $\mu_{1\mu_{-1}}$ is equal to $L$, which only leaves two possibilities for $p$ (and $1 - p$):

$$p = \frac{1}{2} \pm \frac{\rho + 1}{\rho - 1} \sqrt{\frac{1}{4} - \frac{\rho L}{(\rho + 1)^2}}$$

where $\rho = M/m$. Applying the transformation $T$, we get $\nu_{k_{i+1}} = T(\nu_{k_i}) \to T(\nu^*_p) = \nu^*_{1-p}$.

A4. Proof of Theorem 23
(i) It is straightforward to check that $T^2(\nu^*_p) = \nu^*_p$, $\forall p \in (0, 1)$.
(ii) We assume that $SS_A$ is not reduced to $\{m, M\}$ (otherwise $I_u = \emptyset$). We have $\nu_{k+2}(d\lambda) = H(\nu_k, \lambda)\nu_k(d\lambda)$, with

$$H(\nu_k, \lambda) = \frac{(\lambda - \mu_k^p)^2(\lambda - \mu_{k+1}^p)^2}{D_kD_{k+1}} \tag{29}$$

see [15], with $\mu_k^p, D_k$ defined as in Theorem [4]. For $\nu_k = \nu_p^*$, it gives

$$H(\nu_p^*, \lambda) = \frac{[M(1-p) + mp - \lambda]^2[mp + m(1-p) - \lambda]^2}{p^2(1-p)^2(M - m)^4}. \tag{30}$$

One can then check that for any $p \in I_u$, $\max_{\lambda \in SS_A} H(\nu_p^*, \lambda) = H(\nu_p^*, \lambda^*) > 1$, with $\lambda^* = \min_{\lambda \in SS_A} s(\lambda)$. Therefore, for any $p \in I_u$, one can choose $\epsilon$ small enough, such that $d(\nu_k, \nu_p^*) < \epsilon$ implies $\nu_{k+2}([a, b]) > K_p\nu_k([a, b])$, for some $K_p > 1$ and some $a, b$ such that $m + \epsilon < a < b < M - \epsilon$ and $[a, b] \cap SS_A \neq \emptyset$. For any $\alpha > 0$, $\alpha < 1 - p$, take an initial measure $\nu_0$ putting mass $p$ at $m$, $1 - p - \alpha$ at $M$ and $\alpha$ in the interval $[a, b]$. It satisfies $d(\nu_0, \nu_p^*) < \alpha$, and, for any $m$, either $d(\nu_2m, \nu_p^*) < \epsilon$ or $d(\nu_{2(m+1)}([a, b]) > K_p\nu_{2m}([a, b])$. The later case gives $\nu_{2m}([a, b]) > 2\epsilon$, and thus $d(\nu_{2m}, \nu_p^*) < \epsilon$, as soon as $m > \log(2\epsilon/\alpha)/\log(K_p)$, which shows that $\nu_p^*$ is unstable.

(iii) Part (a) concerns the case where a spectral gap is present, with point mass at $m$ and $M$. The proof for the general situation is more technical and is sketched in part (b).

(a) Assume that the measure $\nu_0$ has a spectral gap: $\nu_0 = 0$ on $(m, m+s)$ and $(M - s, M)$ for some $s > 0$. Take $\gamma < s$ and assume that $d(\nu_0, \nu_p^*) < \alpha < \gamma$ with $p \in I_s$. The arguments go as follows. First we bound $\nu_2\{(m+\gamma, M-\gamma)\}$ by $2K_0\alpha$ for some $K_0 < 1$, then we bound $\nu_2\{(M-\gamma, M)\}$ by $1 - p + K_1\alpha$ for some $K_1 < \infty$. We show that $d(\nu_2, \nu_p^*) < K_0\alpha$ for some $p_2$ such that $|p_2 - p| < (K_0 + K_1)\alpha$. Stability will then follow by an induction argument.

The maximum value of $H(\nu_0, \lambda)$ for $\lambda$ varying in $[m + \gamma, M - \gamma]$ may be reached for some $\lambda^* \in (\mu_1^p, \mu_2^p)$ or at one of the two points $m + \gamma, M - \gamma$. Now, for $\alpha$ small enough $H(\nu_0, \lambda)$ will be close to $H(\nu_p^*, \lambda)$ given by (30), and $p \in I_s$ implies

$$\max_{\lambda \in SS_A \cap (\mu_1^p, \mu_2^p)} H(\nu_0, \lambda) < 1. \tag{31}$$

Consider the function $H(\nu_0, \lambda)$ at $\lambda = M - \gamma$. We can write

$$H(\nu_0, M - \gamma) = H(\nu_p^*, M) - \gamma \frac{dH(\nu_p^*, \lambda)}{d\lambda}_{\lambda = M} + F_H(\nu_p^*; \nu_0, M) + \mathcal{O}(\gamma^2), \tag{32}$$

with $F_H(\nu_p^*; \nu_0, M)$ the directional derivative of $H(\nu, M)$ at $\nu_p^*$ in the direction $\nu_0$,

$$F_H(\nu_p^*; \nu_0, M) = \lim_{\beta \to 0^+} \frac{H((1 - \beta)\nu_p^* + \beta\nu_0, M) - H(\nu_p^*, M)}{\beta}. \tag{33}$$
Define $F_H(\nu_p^*; x, \lambda) = F_H(\nu_p^*; \delta_x, \lambda)$ with $\delta_x$ the delta measure supported at $x$. We have

$$F_H(\nu_p^*; \nu_0, M) = \int_m^M F_H(\nu_p^*, x, M) \nu_k(dx),$$

which we decompose in three parts:

$$F_H(\nu_p^*; \nu_0, M) = \int_m^{m+\gamma} F_H(\nu_p^*, x, M) \nu_0(dx) + \int_{m+\gamma}^{M-\gamma} F_H(\nu_p^*, x, M) \nu_0(dx) + \int_{M-\gamma}^M F_H(\nu_p^*, x, M) \nu_0(dx).$$

Direct calculation gives

$$F_H(\nu_p^*, x, M) = \frac{(x-m)^2(M-x)[x-m+(2p-1)(M-m)]}{p^2(1-p)^2(M-m)^4}$$

so that $F_H(\nu_p^*, m, M) = F_H(\nu_p^*, M, M) = 0$ and $F_H(\nu_p^*; \nu_0, M) < F^*\nu_0\{(m+\gamma, M-\gamma)\}$ with $F^* = \max_{x \in \mathbb{L}, \nu \in [m,M]} F_H(\nu_p^*, x, M) < \infty$. Also, $d(\nu_0, \nu_p^*) < \alpha$ implies $\nu_0\{(m+\gamma, M-\gamma)\} = \nu_0\{(m+\alpha, M-\alpha)\} < 2\alpha$, so that $F_H(\nu_p^*; \nu_0, M) < 2\alpha F^*$. Now,

$$H(\nu_p^*, M) = 1, \quad \frac{dH(\nu_p^*, \lambda)}{d\lambda} |_{\lambda=M} = \frac{2}{p(1-p)(M-m)},$$

which, together with (32) gives for $\gamma$ small enough

$$H(\nu_0, M-\gamma) < 1 + 2\alpha F^* - \gamma \frac{1}{p(1-p)(M-m)}$$

and thus

$$H(\nu_0, M-\gamma) < 1 - \frac{\gamma}{2p(1-p)(M-m)}$$

for $\alpha < \gamma/[4F^*p(1-p)(M-m)]$.

The situation is similar at $m+\gamma$. Together with (31) this implies for $\alpha$ small enough

$$\max_{\lambda \in SS, \lambda \in [m+\gamma, M-\gamma]} H(\nu_0, \lambda) < K_0 < 1$$

and therefore

$$\nu_2\{(m+\gamma, M-\gamma)\} < 2K_0 \alpha \quad (33)$$

with $K_0 < 1$ not depending on $\alpha$.

Consider now the interval $(M-\gamma, M]$. We have

$$\nu_2\{(M-\gamma, M]\} = \nu_2(M) = H(\nu_0, M) \nu_0\{(M-\gamma, M]\} = H(\nu_0, M) \nu_0(M),$$

with $d(\nu_0, \nu_p^*) < \alpha$ implying $\nu_0(M) < 1 - p + \alpha$, and

$$H(\nu_0, M) = H(\nu_p^*, M) + F_H(\nu_p^*; \nu_0, M) + \mathcal{O}(\alpha^2) < 1 + 2\alpha F^* + \mathcal{O}(\alpha^2).$$
This gives for $\alpha$ small enough
\[ \nu_2([M - \gamma, M]) < 1 - p + K_1 \alpha \]
for some $K_1 < \infty$. Similarly, $\nu_2([m, m + \gamma]) < p + K_1 \alpha$.

Define $p_0 = p$, $p_2 = [\nu_2(m) - \nu_2(M) + 1]/2$, $\alpha_0 = \alpha$. We obtain
\[-K_0 \alpha_0 < \nu_2(m) - p_2 < 0, \quad -K_0 \alpha_0 < \nu_2(M) - (1 - p_2) < 0\]
which together with \[33\] implies
\[ d(\nu_2, \nu_2^p) < \alpha_2 = K_0 \alpha_0. \]
Moreover, $|p_2 - p_0| < (K_0 + K_1) \alpha_0$.

For $\alpha$ small enough, $p_2 \in \mathcal{I}_s$ and we can then repeat the same arguments. This gives for any $m$
\[ d(\nu_{2m}, \nu_{2m}^p) < \alpha_{2m} = K_0^m \alpha \]
with
\[ |p_{2m} - p| < (K_0 + K_1) \sum_{i=0}^{m-1} \alpha_{2i} = (K_0 + K_1) \frac{1 - K_0^m}{1 - K_0} \alpha < \frac{K_0 + K_1}{1 - K_0} \]
and $p_{2m} \in \mathcal{I}_s$, for $\alpha$ small enough. For any $p \in \mathcal{I}_s$ and any $\epsilon > 0$, taking $\nu_0$ such that $d(\nu_0, \nu_0^p) < \alpha$ with $\alpha$ small enough thus implies $d(\nu_{2m}, \nu_{2m}^p) < \epsilon$ for any $m$, and $\nu_{2m}^p$ is thus stable.

(b) Consider now the general situation. The proof follows the same lines as in case (a), but more technicalities are required since we need to consider measures of intervals of the form $[m, m + \gamma]$ and $(M - \gamma, M]$, with $\gamma$ decreasing in a suitable way as the number of iterations of the mapping $T^2$ increases.

Assume that
\[ \nu_{2k}([m + \gamma_{2k}, M - \gamma_{2k}]) < 2 \alpha_{2k}, \]
\[ \nu_{2k}([m, m + \gamma_{2k}]) < p_{2k} + \alpha_{2k}, \]
\[ \nu_{2k}((M - \gamma_{2k}, M]) < 1 - p_{2k} + \alpha_{2k} \]
for some $p_{2k} \in \mathcal{I}_s$ and some $\alpha_{2k}, \gamma_{2k}$. Note that it implies $d(\nu_{2k}, \nu_{2k}^p) < \gamma_{2k}$ and that for $k = 0$, $\alpha_0, \gamma_0$ can be chosen arbitrarily small, with $d(\nu_0, \nu_0^p) < \alpha_0$ for some $p \in \mathcal{I}_s$.

Consider one application of the mapping $T^2$ at a generic iteration $k$. We can write $H(\nu_{2k}, M) = H(\nu_{2k}^*, M) + F_H(\nu_{2k}^*, \nu_{2k}, M) + O(\gamma_{2k}^2)$ with
\[ F_H(\nu_{2k}^*; \nu_{2k}, M) = \int_0^{m+\gamma_{2k}} F_H(\nu_{2k}^*, x, M)\nu_{2k}(dx) \]
\[ + \int_{m+\gamma_{2k}}^{M-\gamma_{2k}} F_H(\nu_{2k}^*, x, M)\nu_{2k}(dx) \]
\[ + \int_{M-\gamma_{2k}}^{M} F_H(\nu_{2k}^*, x, M)\nu_{2k}(dx). \]
Asymptotic behaviour of a family of gradient algorithms in $\mathbb{R}^d$ and Hilbert spaces

The first integral term is of the order $O(\gamma_{2k}^2)$ (since $F_H(\nu^*_p, m, M) = 0$ and $dF_H(\nu^*_p, z, M)/dz_{|z=m} = 0$), the second is bounded by $2\alpha_2 F^* + O(\gamma_{2k}^2)$, as in case (a). For the third term, for which $x$ is close to $M$, we can use the linear approximation

$$F_H(\nu^*_p, x, M) = (x - M)\frac{dF_H(\nu^*_p, z, M)}{dz} \bigg|_{z=M} + O(\gamma_{2k}^2)$$

which gives

$$\int_{M-\gamma_{2k}}^{M} F_H(\nu^*_p, x, M)\nu_2k(dx) = \frac{2}{p_2k(1-p_2k)^2(M-m)} I_2k(M) + O(\gamma_{2k}^2)$$

where $I_2k(M) = \int_{0}^{\gamma_{2k}} z\nu'_2k(dz)$ with $\nu'_2k$ the measure obtained after applying the transformation $x \rightarrow z = M - x$. We have thus obtained

$$H(\nu_2k, M) < 1 + 2\alpha_2 F^* + \frac{2I_2k(M)}{p_2k(1-p_2k)^2(M-m)} + O(\gamma_{2k}^2). \quad (34)$$

Consider now the behavior of $I_2k(M)$ as $k$ increases. We assume that $\nu_2k$ remains in some neighborhood $\mathcal{V}(p)$ of $\nu^*_p$, which we shall be able to guarantee afterwards. Define $A_{2k}(M) = I_2k(M)[\int_{0}^{\gamma_{2k}} \nu'_2k(dz)]^{-1}$. It satisfies $I_2k(M) < A_{2k}(M) < \gamma_{2k}$. Also, $\gamma_{2(k+1)} < \gamma_{2k}$ implies

$$A_{2(k+1)}(M) = \frac{\int_{0}^{\gamma_{2(k+1)}} zH(\nu_2k, M - z)\nu'_2k(dz)}{\int_{0}^{\gamma_{2(k+1)}} H(\nu_2k, M - z)\nu'_2k(dz)} < \frac{\int_{0}^{\gamma_{2k}} zH(\nu_2k, M - z)\nu'_2k(dz)}{\int_{0}^{\gamma_{2k}} H(\nu_2k, M - z)\nu'_2k(dz)},$$

and, since $H(\nu_2k, M - z)$ decreases for $z$ close to zero,

$$A_{2(k+1)}(M) < \frac{\int_{0}^{\gamma_{2k}} z(1 - az)\nu'_2k(dz)}{\int_{0}^{\gamma_{2k}} \nu'_2k(dz)}.$$

We can bound the speed of decrease of $H(\nu_2k, M - z)$: $H(\nu, M - z)/H(\nu, M) < 1 - az$ for some $a > 0$, any $z$ in $[0, \gamma_0]$ and any $\nu \in \mathcal{V}(p)$. This gives

$$A_{2(k+1)}(M) < \frac{\int_{0}^{\gamma_{2k}} (1 - az)\nu'_2k(dz)}{\int_{0}^{\gamma_{2k}} \nu'_2k(dz)}.$$

Repeating the same arguments we get for any $n > 0$,

$$A_{2(k+n)}(M) < \tilde{A}_{2(k+n)}(M) = \frac{\int_{0}^{\gamma_{2k}} (1 - az)^n\nu'_2k(dz)}{\int_{0}^{\gamma_{2k}} \nu'_2k(dz)},$$

with $\tilde{A}_{2(k+n)}(M)$ decreasing with $n$. Direct calculation gives $\sum_{n=0}^{\infty} \tilde{A}_{2(k+n)}(M) = 1/a$, and therefore $I_{2k}(M) < \tilde{A}_{2k}(M) = o(1/k)$.
Similarly to case (a), we can write

\[ H(\nu_{2k}, M - \gamma_{2(k+1)}) = H(\nu_{2k}, M) - \frac{2\gamma_{2(k+1)}}{p_{2k}(1 - p_{2k})(M - m)} + O(\gamma_{2k}^2), \]

with \( H(\nu_{2k}, M) \) bounded by (34). Assume that \( \gamma_{2k} \) is such that \( A_{2k}(M) = o(\gamma_{2k}) \) and \( \alpha_{2k} = o(\gamma_{2k}) \). We obtain for \( p_{2k} \) close enough to \( p \)

\[ H(\nu_{2k}, M - \gamma_{2(k+1)}) < \beta_{2(k+1)} = 1 - \frac{\gamma_{2(k+1)}}{p(1 - p)(M - m)}. \]  

(35)

We thus get the following bounds on the measure of subintervals of interest at the next iteration:

\[ \nu_{2(k+1)}\{ (m + \gamma_{2(k+1)}, M - \gamma_{2(k+1)}) \} < 2 \max\{ \beta_{2(k+1)}, K_0 \} \alpha_{2k} \]  

(36)

where \( K_0 = \max_{p_{2k} \in \nu(p)} \max_{\lambda \in S \cap \{ (\mu^2_{2k}, \nu_{2k}) \}} H(\nu_{2k}, \lambda) \), and \( K_0 < 1 \) for \( p \) in \( \mathcal{I}_s \) and \( \mathcal{V}(p) \) small enough, see part (a);

\[ \nu_{2(k+1)}\{ (M - \gamma_{2(k+1)}, M) \} < \nu_{2(k+1)}\{ (M - \gamma_{2k}, M) \} \]

\[ < H(\nu_{2k}, M)\nu_{2k}\{ (M - \gamma_{2k}, M) \} \]

\[ < \left[ 1 + 2\alpha_{2k} F^* + \frac{2A_{2k}(M)}{p_{2k}(1 - p_{2k})^2(M - m)} + O(\gamma_{2k}^2) \right] \nu_{2k}\{ (M - \gamma_{2k}, M) \} \]

\[ < \nu_{2k}\{ (M - \gamma_{2k}, M) \} + B\alpha_{2k} + CA_{2k}(M) + D\gamma_{2k}^2 \]

for some \( B, C, D < \infty \). Similarly, we obtain

\[ \nu_{2(k+1)}\{ (m, m + \gamma_{2(k+1)}) \} < \nu_{2k}\{ (m, m + \gamma_{2k}) \} + B\alpha_{2k} + CA_{2k}(m) + D\gamma_{2k}^2 \]

where \( A_{2k}(m) \) is defined similarly to \( A_{2k}(M) \). Define \( p_{2(k+1)} \) as

\[ p_{2(k+1)} = \frac{\nu_{2(k+1)}\{ (m, m + \gamma_{2(k+1)}) \} - \nu_{2(k+1)}\{ (M - \gamma_{2(k+1)}, M) \} + 1}{2}, \]

it gives

\[ 0 < p_{2(k+1)} - \nu_{2(k+1)}\{ (m, m + \gamma_{2(k+1)}) \} < \max\{ \beta_{2(k+1)}, K_0 \} \alpha_{2k}, \]

\[ 0 < 1 - p_{2(k+1)} - \nu_{2(k+1)}\{ (M - \gamma_{2(k+1)}, M) \} < \max\{ \beta_{2(k+1)}, K_0 \} \alpha_{2k}. \]

Together with (33) it implies

\[ \left| p_{2(k+1)} - p_{2k} \right| < \Delta_{2k} = [B + 1 + \max\{ \beta_{2(k+1)}, K_0 \}] \alpha_{2k} + CA'_{2k} + D\gamma_{2k}^2, \]

where \( A'_{2k} = \max\{ A_{2k}(m), A_{2k}(M) \} \) and \( \sum\Delta_{2k} < \infty \).

Define \( \alpha_{2(k+1)} = \max\{ \beta_{2(k+1)}, K_0 \} \alpha_{2k} \) and take \( \gamma_{2k} = 1/k^q \) with \( q < 1 \), so that \( A'_{2k} = o(\gamma_{2k}) \). From the definition of \( \beta_{2(k+1)} \), see (35), \( \sum\kappa \alpha_{2k} < \infty \) and \( \alpha_{2k} = o(\gamma_{2k}) \). Since \( \sum\kappa A_{2k}^2 < \infty \), taking \( q > 1/2 \) in the definition of \( \gamma_{2k} \) ensures \( \sum\Delta_{2k} < \infty \). We can repeat the same argument, and \( d(\nu_{2(k+n)}, \nu_{p_{2(k+n)}}) < \gamma_{2(k+n)} \) which tends to zero as \( n \) increases, with \( \left| p_{2(k+n)} - p_{2k} \right| \) remaining finite.
\( \nu_2(k+n) \) thus remains in some neighborhood \( \mathcal{V}(p) \) of \( \nu_1^* \) for any \( n \), and \( \mathcal{V}(p) \) can be made arbitrarily small by choosing \( \alpha_0 \) and \( \gamma_0 \) small enough. \hfill \blacksquare

**A5. Proof of Theorem 6.** Assume that \( x_0 \) is such that for some \( k \geq 0 \), \( \|g_{k+1}\| = 0 \) with \( \|g_i\| > 0 \) for all \( i \leq k \) (that is, \( x_{k+1} = x^* \) and \( x_i \neq x^* \) for \( i \leq k \)). This implies \( R_k(W) = 0 \) for any \( W \), and therefore \( R(W, x_0, x^*) = R(x_0, x^*) = 0 \).

Assume now that \( \|g_k\| > 0 \) for all \( k \). Consider

\[
V_n = \left[ \prod_{k=0}^{n-1} R_k(W) \right]^{1/n} = \left[ \prod_{k=0}^{n-1} \frac{(Wg_{k+1}, g_{k+1})}{(Wg_k, g_k)} \right]^{1/n} = \left( \frac{(Wg_n, g_n)}{(Wg_0, g_0)} \right)^{1/n}.
\]

We have,

\[
\forall z \in \mathcal{H}, \ c\|z\|^2 \leq (Wz, z) \leq C\|z\|^2,
\]

and thus

\[
(c/C)^{1/n} \left( \frac{g_n}{g_0} \right)^{1/n} \leq V_n \leq (C/c)^{1/n} \left( \frac{g_n}{g_0} \right)^{1/n}.
\]

Since \((c/C)^{1/n} \to 1\) and \((C/c)^{1/n} \to 1\) as \( n \to \infty \), \( \lim \inf_{n \to \infty} V_n \) and \( \lim \sup_{n \to \infty} V_n \) do not depend on \( W \). Take \( W = P(A) \); it gives \( R_k(W) = r_k = 1 - 1/L_k \), see [9], which is not decreasing, and thus \( \lim_{n \to \infty} V_n = 1 - 1/L \) for any \( W \). \hfill \blacksquare

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