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Geometric Numerical Methods and Results in the Contrast Imaging Problem in Nuclear Magnetic Resonance

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The purpose of this article is to present numerical methods and results about the contrast imaging problem in nuclear magnetic resonance which corresponds to a Mayer problem in optimal control. The candidates as minimizers are selected among a set of extremals, solutions of an Hamiltonian system given by the Pontryagin Maximum Principle and sufficient second order conditions are described. They form the geometric foundations of the HAMPATH code which combines shooting and continuation methods, see Ref. 12. The main contribution of this article is to present a numerical analysis of the contrast imaging problem in NMR in the case of deoxygenated/oxygenated blood samples as an application of the aforementioned techniques.

Keywords: Geometric optimal control; Contrast imaging in NMR; Shooting and continuation methods.

AMS Subject Classification: 49K15, 49M15, 81V55

1. Introduction

Geometric optimal control was introduced very recently in nuclear magnetic resonance (NMR) to control the dynamics of spin particles, e.g. see Ref. 10. The model is the coupling of a system equations introduced by F. Bloch in 1946, describing

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2 B. BONNARD, O. COTS

the motion of nuclear magnetization:

$$\begin{aligned}
 \frac{dM_x}{dt} &= -\frac{M_x}{T_2} - \Delta\omega M_y + \omega_y M_z \\
 \frac{dM_y}{dt} &= -\frac{M_y}{T_2} + \Delta\omega M_x - \omega_x M_z \\
 \frac{dM_z}{dt} &= \frac{(M_0 - M_z)}{T_1} - \omega_y M_x + \omega_x M_y,
 \end{aligned} \tag{1.1}$$

where $M = (M_x, M_y, M_z)$ is the magnetization vector, M_0 represents the equilibrium state of the system, T_1, T_2 are the *relaxation times* characteristic of the particle, $\Delta\omega$ is the *detuning* and $\omega = (\omega_x, \omega_y)$ is the control which is a transverse radio frequency field in the (x, y) plane. This model represents the physical experiment in NMR spectroscopy very accurately and the control can be realized in practice as shown, for instance, by the *saturation problem* of a spin 1/2 particle which consists in bringing the magnetization vector to zero, see Ref. 21.

An ideal model in the contrast imaging problem in NMR consists of considering a system of two Bloch Eq. (1.1) with zero detuning and distinct relaxation parameters representing two different spin 1/2 particles, controlled by the same magnetic field and bringing the magnetization of the first particle to zero while maximizing the magnitude of the magnetization vector of the second particle.

This is illustrated by the experimental results, presented on Fig. 1. The sample consists of two test tubes with outer diameters of 5mm and 8mm. The outer and inner volumes were respectively filled with two solutions of oxygenated and deoxygenated blood. The left hand side picture represents the two samples at the equilibrium (they both appear white) and the right hand side picture shows the result after the application of the control sequence. The inner sample appears black (zero magnetization) while the outer sample represents the remaining magnetization. For the details see Ref. 22.



Fig. 1. Experimental results: The inner circle shape sample mimics the deoxygenated blood, where $T_1 = 1.3s$ and $T_2 = 50ms$; the outside moon shape sample corresponds to the oxygenated blood, where $T_1 = 1.3s$ and $T_2 = 200ms$.

In particular, using the symmetry of revolution of each spin system with respect to the z -axis, we shall concentrate to the case where $\omega_y \equiv 0$ and each spin system can be restricted to the (y, z) plane, defining the system:

$$\begin{aligned}\frac{dM_y^i}{dt} &= -\frac{M_y^i}{T_2^i} - \omega_x M_z^i \\ \frac{dM_z^i}{dt} &= \frac{(M_0^i - M_z^i)}{T_1^i} + \omega_x M_y^i, \quad i = 1, 2.\end{aligned}\tag{1.2}$$

The contrast problem is then the following optimal control problem: transfer the system from the equilibrium state in a given transfer time t_f to $M^1(t_f) = 0$ while maximizing $|M^2(t_f)|^2$ and $|M^2(t_f)|$ represents the contrast.

This problem corresponds to a *Mayer problem* in optimal control:

$$\min_{u(\cdot)} c(x(t_f)), \quad \frac{dx}{dt} = f(x, u), \quad u \in U,$$

with the boundary conditions:

$$x(0) = x_0, \quad g(x(t_f)) = 0$$

and it can be analyzed using Pontryagin Maximum Principle, see Ref. 4. It tells us that optimal solutions are found as projections on the x -space as *extremals* solutions of the Hamiltonian system:

$$\begin{aligned}\frac{dx}{dt} &= \frac{\partial H}{\partial p}(x, p, u), \quad \frac{dp}{dt} = -\frac{\partial H}{\partial x}(x, p, u) \\ H(x, p, u) &= \max_{v \in U} H(x, p, v),\end{aligned}\tag{1.3}$$

where $H(x, p, u) = \langle p, f(x, u) \rangle$ is the *pseudo-Hamiltonian* and $p(\cdot)$ is the adjoint vector. Moreover the solution has to satisfy the *transversality condition*:

$$p(t_f) = p^0 \frac{\partial c}{\partial x}(x(t_f)) + \sum_{i=1}^k \sigma_i \frac{\partial g_i}{\partial x}(x(t_f)), \quad p^0 \leq 0.$$

The Maximum Principle is only a necessarily optimality condition and it was recently completed by generalizing the concept of *conjugate point* and *extremal field* in standard calculus of variation, see Ref. 16, to get sufficient geometric optimality conditions, see Ref. 7.

All this theoretical framework will be rapidly recalled in this article before presenting the main contribution of this article consisting in describing the algorithms and simulations in one specific example in the contrast imaging problem. The general method which is the core of the HAMPATH code Ref. 13 is the following.

The Maximum Principle together with the transversality condition leads to the computation of the optimal solution using a *shooting equation*: $S(p(0)) = 0$, integrating the ODE according to Eq. (1.3) and solving with prescribed boundary conditions to compute an initial condition on the adjoint vector $p(0)$. If the method is simple, an initial guess has to be determined to ensure the convergence of the

4 B. BONNARD, O. COTS

shooting. Hence the shooting is complemented by a *smooth continuation method*. The Hamiltonian dynamics defined by Eq. (1.3) is embedded into a one-parameter family of Hamiltonian dynamics H_λ , $\lambda \in [0, 1]$ and this leads to a one-parameter family of shooting equations: $S(p(0), \lambda) = 0$. The connected component of the set of zeros $p_\lambda(0)$ will form a differential curve which is followed from $\lambda = 0$ to $\lambda = 1$ to compute the solution. Also the differential of this curve is shown to be related to the concept of conjugate point.

In order to be applied to the contrast imaging problem one must adapt this general algorithm. Indeed in the general case of Eq. (1.1) due to the symmetry of revolution there is a one parameter family of solutions of the shooting equation. One adaptation concerning the case treated in this article restricting the control to $\omega_y \equiv 0$ is the following. This case leads to a *single-input* control system of the form:

$$\frac{dx}{dt} = F(x) + uG(x), \quad |u| \leq 1.$$

In the single-input case, the optimal solution is a concatenation of *regular arcs* where the control is given by $u(t) = \text{sgn}\langle p(t), G(x(t)) \rangle$ and *singular arcs* where $\langle p(t), G(x(t)) \rangle = 0$, identically. An optimal solution is a concatenation of bang arcs with $u = \pm 1$ and singular arcs and the problem is to determine the sequence of such arcs depending upon the transfer time and the relaxation parameters.

For that the technique is to use a regularization of the optimal control problem transforming the Mayer problem into a *Bolza problem*, for instance with cost:

$$c(x(t_f)) + (1 - \lambda) \int_0^{t_f} |u|^{2-\lambda} dt, \quad \lambda \in [0, 1].$$

Once the structure of the solution is known, it is computed using a multiple-shooting method. In particular, in the case of a deoxygenated/oxygenated blood samples, one proves in this article that the optimal solution is the concatenation of a bang and a singular arc which is finally accurately computed.

The organization of this article is the following. In section 2, we present in details the model used in the contrast imaging problem. In section 3, we recall the necessary and sufficient optimality conditions with applications to our problem. The main results in this article are finally presented in sections 4 and 5, which respectively describe the algorithmic schemes and the numerical simulations in the study case of deoxygenated/oxygenated blood samples.

2. The mathematical model and the Mayer optimal control problem

According to Eq. (1.2) using normalized coordinates, one considers the system formed by coupling two spin 1/2 systems with zero detuning, described for $i = 1, 2$

by:

$$\begin{aligned} \frac{dy_i}{dt} &= -\Gamma_i y_i - uz_i \\ \frac{dz_i}{dt} &= \gamma_i(1 - z_i) + uy_i, \end{aligned} \tag{2.1}$$

with $\Gamma_i = 2\pi/(\omega_{\max}T_2^i)$, $\gamma_i = 2\pi/(\omega_{\max}T_1^i)$ while the control is bounded by $|u| \leq 2\pi$. From the experimental point of view, $\omega_{\max}/2\pi$ can be chosen up to 15000 Hz but the value 32.3 Hz will be considered in our experiments, see Ref. 21 for more details. For the experiments one considers four cases, the relaxation times being:

- | | | |
|----------------------------|--------------------|-----------------|
| 1. Cerebrospinal fluid | : $T_1 = 2000$ ms, | $T_2 = 200$ ms |
| Water | : $T_1 = 2500$ ms, | $T_2 = 2500$ ms |
| 2. Deoxygenated blood | : $T_1 = 1350$ ms, | $T_2 = 50$ ms |
| Oxygenated blood | : $T_1 = 1350$ ms, | $T_2 = 200$ ms |
| 3. Gray matter of cerebrum | : $T_1 = 920$ ms, | $T_2 = 100$ ms |
| White matter of cerebrum | : $T_1 = 780$ ms, | $T_2 = 90$ ms |
| 4. Water | : $T_1 = 2500$ ms, | $T_2 = 2500$ ms |
| Fat | : $T_1 = 200$ ms, | $T_2 = 100$ ms |

Denoting $x_1 = (y_1, z_1)$, $x_2 = (y_2, z_2)$ as the vectors representing each spin, the Bloch ball $|x_i| \leq 1$ being invariant for the dynamics since the parameters satisfy $2\Gamma_i \geq \gamma_i$, the optimal control problem is the following: starting from the north pole $N = ((0, 1), (0, 1))$ corresponding to the equilibrium point of the uncontrolled dynamics, the goal is to reach in a given transfer time t_f the final state $x_1(t_f) = 0$ corresponding to zero magnetization of the first spin, while maximizing the square norm $|x_2(t_f)|^2$ of the second spin.

Introducing the system $\dot{x} = f(x, u)$, where $x = (x_1, x_2)$ belongs to $|x_i| \leq 1$ and $|u| \leq 2\pi$, formed by the coupling of two equations Eq. (2.1), with $x(0) = N$, $g(x(t_f)) = x_1(t_f)$ and with the cost functions $c(x(t_f)) = -|x_2(t_f)|^2$, the problem is a standard Mayer problem given by:

- a system $\dot{x} = f(x, u)$, $u \in U$, $x(0) = x_0$, where U is the control domain and \mathcal{U} is the set of admissible controls,
- a final boundary condition $g(x(t_f)) = 0$, and
- a cost to minimize, $\min_{u(\cdot) \in \mathcal{U}} c(x(t_f))$.

6 *B. BONNARD, O. COTS*

3. Necessary and sufficient optimality conditions for a Mayer problem

3.1. Notation

Let X be an open manifold of \mathbb{R}^n , $T_x X$, $T_x^* X$ the tangent and cotangent spaces. If F, G are two smooth vector fields, the Lie bracket is given by:

$$[F, G] = \frac{\partial F}{\partial x}(x)G(x) - \frac{\partial G}{\partial x}(x)F(x),$$

while the Poisson bracket of two smooth Hamiltonians H_1, H_2 (the associated Hamiltonian vector fields being denoted \vec{H}_i) is defined by $\{H_1, H_2\}(z) = dH_1(\vec{H}_2)$, $z = (x, p)$. Denoting H_F, H_G the Hamiltonian lifts of F, G : $H_F = \langle p, F(x) \rangle$, $H_G = \langle p, G(x) \rangle$, one has the relation $\{H_F, H_G\} = \langle p, [F, G](x) \rangle$.

3.2. Geometric preliminaries

To highlight the necessary optimality conditions, one introduces the following geometric framework. Consider a (smooth) Mayer problem of the form:

$$\min_{u(\cdot) \in \mathcal{U}} c(x(t_f)), \quad \frac{dx}{dt} = F(x) + uG(x),$$

$x \in X \subset \mathbb{R}^n$, $|u| \leq 1$, with fixed initial condition: $x(0) = x_0$ with the terminal condition $x(t_f) \in M$, where M is a submanifold of codimension k of X defined by: $g(x) = 0$. The class of admissible controls \mathcal{U} is the set of L^∞ -mappings, valued in $U : |u| \leq 1$ such that the solution $x(\cdot, x_0, u)$ is defined on $[0, t_f]$. For $t \leq t_f$ one defines the accessibility set at time t : $A(x_0, t) = \bigcup_{u(\cdot) \in \mathcal{U}} x(t, x_0, u)$. Assuming that the cost c is a regular mapping, such that for $c(x) = m$, the set $M_m = \{x; g(x) = 0, c(x) = m\}$ is a submanifold of codimension $k + 1$, one gets the following necessary optimality condition according to the Pontryagin Maximum Principle and its extension, see Refs. 19 and 4.

3.3. Maximum Principle and High Order Maximum Principle

Theorem 3.1. *Let $u^*(\cdot)$ be an admissible control whose corresponding trajectory $x^*(\cdot)$ is optimal on $[0, t_f]$. Then there exists a non-zero adjoint vector function $p^*(\cdot)$ such that the conditions*

$$\frac{dx^*}{dt} = \frac{\partial H}{\partial p}(x^*, p^*, u^*), \quad \frac{dp^*}{dt} = -\frac{\partial H}{\partial x}(x^*, p^*, u^*) \quad (3.1)$$

are satisfied almost everywhere, where $H = \langle p, F + uG \rangle$ is the pseudo-Hamiltonian. Denoting $M(x, p) = \max_{|u| \leq 1} H(x, p, u)$ one must have a.e. the maximization condition

$$H(x^*, p^*, u^*) = M(x^*, p^*) \quad (3.2)$$

and moreover M is a constant.

In the singular case where $H_G = 0$ and if the control is not saturating, i.e. $|u^*| < 1$, the generalized Legendre–Clebsch condition,

$$\{H_G, \{H_G, H_F\}\} \leq 0 \quad (3.3)$$

has to be satisfied.

Finally we have the boundary conditions

$$g(x^*(t_f)) = 0 \quad (3.4)$$

and

$$p^*(t_f) = p_0 \frac{\partial c}{\partial x}(x^*(t_f)) + \sum_{i=1}^k \sigma_i \frac{\partial g_i}{\partial x}(x^*(t_f)), \quad (3.5)$$

$p_0 \leq 0$, $\sigma = (\sigma_1, \dots, \sigma_k) \in \mathbb{R}^k$. This last condition is called the transversality condition.

Definition 3.1. We call any solution of Eq. (3.1) and Eq. (3.2) an extremal, and if the boundary conditions Eqs. (3.4), (3.5) are satisfied they are called BC-extremals. Condition (3.2) imposes $u(t) = \text{sgn}\langle p(t), G(x(t)) \rangle$ at any time such that $\langle p(t), G(x(t)) \rangle \neq 0$. An extremal is called regular if $u(t) = \text{sgn} H_G$ and bang-bang if the number of switchings is finite. It is called singular if $H_G = 0$.

3.4. Computation of singular extremals

This computation is standard: let $z(\cdot) = (x(\cdot), p(\cdot))$ be an extremal, $t \in [0, t_f]$. In the singular case, we have $H_G(z(t)) = 0$ identically. Differentiating with respect to time,

$$H_G(z(t)) = \{H_G, H_F\}(z(t)) = 0 \quad (3.6)$$

$$\{\{H_G, H_F\}, H_F\}(z(t)) + u_s(t)\{\{H_G, H_F\}, H_G\}(z(t)) = 0 \quad (3.7)$$

and Eq. (3.7) defines the singular control u_s as a dynamic feedback, outside $S = \{z; \{\{H_G, H_F\}, H_G\}(z) = 0\}$. Denote Σ the switching surface $H_G = 0$ and $\Sigma' : H_G = \{H_G, H_F\} = 0$. Plugging u_s into H defines a true Hamiltonian function denoted H_s outside S . We have:

Proposition 3.1. *Outside S , the singular extremals are the solutions of*

$$\frac{dx}{dt} = \frac{\partial H_s}{\partial p}, \quad \frac{dp}{dt} = -\frac{\partial H_s}{\partial x},$$

starting at $t = 0$ from $\Sigma' \setminus S$ and they are contained in Σ' .

8 *B. BONNARD, O. COTS*

For the computations of the singular trajectories, we need the following Lie brackets:

$$\begin{aligned}
 F &= \sum_{i=1,2} (-\Gamma_i y_i) \frac{\partial}{\partial y_i} + (\gamma_i(1 - z_i)) \frac{\partial}{\partial z_i} \\
 G &= \sum_{i=1,2} -z_i \frac{\partial}{\partial y_i} + y_i \frac{\partial}{\partial z_i} \\
 [G, F] &= \sum_{i=1,2} (\gamma_i - \delta_i z_i) \frac{\partial}{\partial y_i} - (\delta_i y_i) \frac{\partial}{\partial z_i} \\
 [[G, F], F] &= \sum_{i=1,2} (\gamma_i(\gamma_i - 2\Gamma_i) - \delta_i^2 z_i) \frac{\partial}{\partial y_i} + \delta_i^2 y_i \frac{\partial}{\partial z_i} \\
 [[G, F], G] &= \sum_{i=1,2} 2\delta_i y_i \frac{\partial}{\partial y_i} + (\gamma_i - 2\delta_i z_i) \frac{\partial}{\partial z_i},
 \end{aligned}$$

where $\delta_i = \gamma_i - \Gamma_i$.

3.5. Application to the contrast problem: Boundary conditions

Since the Bloch ball is invariant for the dynamics, one can take $X = \{|x_i| < 1; i = 1, 2\}$. Applying the Maximum Principle, the boundary conditions Eqs. (3.4) and (3.5) give us:

$$\begin{aligned}
 x_1(t_f) &= 0 \quad (\text{zero magnetization of the first spin}) \\
 p_2(t_f) &= -2p^0 x_2(t_f), \quad p^0 \leq 0.
 \end{aligned} \tag{3.8}$$

If p^0 is non-zero, it can be normalized by homogeneity to $p^0 = -1/2$. The case $p^0 = 0$ corresponds to the case where t_f is the minimum time to transfer the first spin to zero.

3.6. Second-order optimality conditions

This section relies to the work of Ref. 7.

Definition 3.2. Let $z(\cdot)$ be a reference singular solution of \vec{H}_s on $[0, t_f]$ and contained in Σ' . The variational equation

$$\begin{aligned}
 \delta \dot{z} &= d\vec{H}_s(z(t))\delta z \\
 dH_G &= d\{H_F, H_G\} = 0
 \end{aligned}$$

is called Jacobi equation. A Jacobi field $J(t) = (\delta x, \delta p)$ is a non-zero solution of the Jacobi equation. It is said semi-vertical at time t if $\delta x(t) \in \mathbb{R}G(x(t))$. The times $0 \leq t_1 < t_2 \leq t_f$ are said to be conjugate if there exists a Jacobi field $J(t)$ semi-vertical at $t = t_1, t_2$ and the points $x_1 = x(t_1)$ and $x_2 = x(t_2)$ are said to be conjugate.

Assumptions Let $z(\cdot) = (x(\cdot), p(\cdot))$ be a reference singular extremal curve on $[0, t_f]$ solution on $\Sigma' \setminus S$ and we assume the following:

- (A1) $t \mapsto x(t)$ is a one-to-one immersion. Applying a feedback, one can identify the reference singular control to zero.
- (A2) *Strong corank one condition.*
 $K(t) = \text{span}\{\text{ad}^k F.G(x(t)); k \geq 0, t \in [0, t_f]\}$ is of codimension one and generated by the first $(n - 1)$ vectors $\{\text{ad}^k F.G(x(t)); k = 0, \dots, n - 2\}$ where $\text{ad} F.G$ denotes $[F, G]$.
- (A3) *Non-exceptional case.* $H_s(z(t))$ is non-zero.

3.6.1. Optimality results

One considers a Mayer problem for a system $\dot{x} = F(x) + uG(x)$, $x \in X$ open set of \mathbb{R}^n , $|u| \leq 2\pi$, $\min_{u(\cdot)} c(x(t_f))$ and one assumes that the terminal manifold M is a trajectory of the vector field G : $\{\exp tG; t \in \mathbb{R}\}$. In this case one will derive a sufficient optimality condition, which is applicable to the contrast imaging problem.

First one recalls the following proposition from Ref. 7.

Proposition 3.2. *Assuming $u \in \mathbb{R}$, a necessary optimality condition in the Mayer problem for a singular arc δ_s on $[0, t_f]$ satisfying (A1), (A2) and (A3) is the non-existence of a time $0 < t_c < t_f$ conjugate to zero, for the solutions of the system contained in a tubular neighbourhood of δ_s .*

Proof. According to Ref. 7, the first conjugate time $0 < t_{1c}$ is the first time t such that the extremity mapping $E^{x_1, t} : u(\cdot) \mapsto x(t, x_1, u)$ becomes open when restricting to curves contained in a tubular neighbourhood of δ_s . This is clearly a necessary optimality condition. \square

Next, one presents the following sufficient optimality condition.

Theorem 3.2. *Consider the Mayer problem with the terminal manifold $M = \{\exp tG; t \in \mathbb{R}\}$. Let $\delta_+ \delta_s$ be a BC-extremal on $[0, t_f]$ formed by a bang arc with $u = 2\pi$ on $[0, t_1]$ followed by a singular arc on $[t_1, t_f]$. Assume that the singular arc $t \mapsto \delta_s(t)$ satisfies the assumptions (A1), (A2), (A3) and is associated to a control $u_s(\cdot)$ on $[t_1, t_f]$ which is not saturating, i.e., $|u_s| < 2\pi$. Then the extremal $\delta_+ \delta_s$ is optimal on $[0, t_f]$ in a tubular neighbourhood of δ_s provided there exists no conjugate point on $[t_1, t_f]$.*

Proof. Denote x_f the terminal point $\delta_s(t_f)$ and p_f the adjoint vector. One has $\langle p_f, G(x_f) \rangle = 0$. According to the classification of extremals due to Ref. 20, near (x_f, p_f) every extremal is of the form $\delta_\pm \delta_s \delta_\pm$, where an arc of the sequence can be empty. Let $M_\varepsilon = \{\exp sG(x_f); |s| < \varepsilon \text{ small enough}\}$. Using the transversality condition, near (x_f, p_f) every BC-extremal starting from M_ε is of the form $\delta_\pm \delta_s$.

10 *B. BONNARD, O. COTS*

Let $S(x_f, t)$, $t \in [t_1, t_f]$ be the surface formed by integrating backwards in time the singular flow initiating from M_ε . For fixed t , it is a surface of codimension two, whose tangent space can be easily computed at a point $\delta_s(t)$, $t \in [t_1, t_f]$. It is a plane $E(t)$ spanned by $d\Pi(J(t))$, $J(t)$ being the Jacobi fields which are semi-vertical at t_f , where Π is the standard projection: $(x, p) \mapsto x$.

Hence if $\text{rank} \{E, F, G\}_{|\delta_s(t)}$ is four for $t < t_f$, then a policy $\delta_\pm \delta_s$ will form an extremal field along the reference singular extremal δ_s . Otherwise, if it is not of full rank, at a time t , then by definition $\delta_s(t)$ and $\delta_s(t_f)$ are conjugate. Now clearly, since conjugate times correspond to points where the extremity mapping becomes open, if $\delta_s(t)$ and $\delta_s(t_f)$ are conjugate for $t_1 \leq t < t_f$, then there exists a pair of conjugate points on $[t_1, t_f]$ and a point conjugate to $x(t_1)$. \square

3.6.2. Application to the contrast problem

First of all in the contrast problem where the control bound is normalized to 2π , one has $\Gamma_i, \gamma_i \ll 1$. Hence a bang solution amounts roughly to a rotation of each plane (y_i, z_i) around 0. Moreover the north pole $N = ((0, 1), (0, 1))$, which is the initial state of the Mayer problem, is an equilibrium point for the singular control system and due to symmetry of revolution, one can take for the first bang either $u = +2\pi$ or $u = -2\pi$. At the final state the first spin has to be at zero, which is the center of the rotation of the first spin. Hence the terminal manifold is indeed a trajectory of the vector field G . In particular, one deduces:

Lemma 3.1. *The simplest BC-extremal in the contrast problem is of the form $\delta_+ \delta_s$.*

3.6.3. The limit case $t_f = t_{min}$, minimum time to transfer the first spin to zero

An interesting situation studied in Ref. 21 is the limit case when the transfer time t_f is exactly the minimum time t_{min} to transfer the first spin to zero. We recall the optimal policy in the physical interesting situation.

Proposition 3.3. *Assume $2\Gamma_1 > 3\gamma_1$, then the time minimal solution to transfer the first spin from the north pole to the center of Bloch ball is of the form $\delta_+ \delta_s \delta_+ \delta_s$ and contains two non-empty singular arcs, the first one being horizontal in the plane (y_1, z_1) while the second is a subarc of the vertical axis of revolution Oz_1 .*

It shows that in the physical application if the transfer time is t_{min} , then the policy $\delta_+ \delta_s$ is not optimal. Also it gives in each experiment an indication about the relation between the contrast and the relaxation parameters. The corresponding contrast is called the Sat-contrast, see Ref. 9. We represent on Fig. 2, the time minimal synthesis to steer the system from the north pole to any point of the Bloch ball. In particular it shows the time minimal solution of the saturation problem.

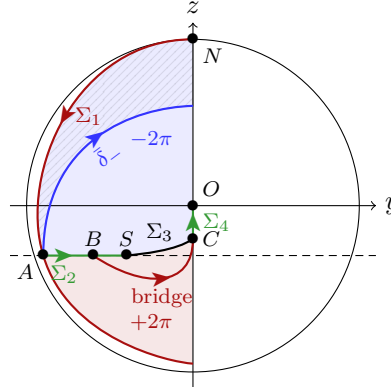


Fig. 2. Schematic time minimal synthesis to steer a single spin system from the north pole N to any point of the Bloch ball. An arbitrary zoom has been used to construct the figure. The set of Σ_i forms the switching surface Σ dividing the $+2\pi$ and -2π areas respectively in red and blue. The δ_- bang arc splits two domains. The hatched one has bang-bang solutions whereas the optimal solutions of the other one contain at least one non trivial singular arc. The minimal time trajectory to steer the spin from N to O (the center of the Bloch ball) is $NABCO$, *i.e.* it is of the form $\delta_+ \delta_s \delta_+ \delta_s$ with vertical and horizontal singular arcs. The spin leaves the horizontal singular arc before the point S (where the control saturates the constraint) producing a bridge to reach the vertical singular line.

4. The numerical methods

In this section we present the numerical methods relating to the resolution of the contrast problem. Each method presented hereunder is a part of the package HAMPATH and all the computations of derivatives are automatically done by the software. The geometrical study of the contrast problem ensures that the optimal solution is a concatenation of bang-singular (BS) sequences. The strategy presented in the following part is first to regularize the problem in order to detect the BS-sequences and find a good approximation of the solution. This approximation obtained by smooth continuation technique is then used to initialize the multiple-shooting computation of the BS-sequences. Up to this point, the transfer duration t_f is fixed to $1.1t_{min}$, where t_{min} is the minimal time to steer the first spin from the north pole to the center of the Bloch ball. Finally, we study the behavior of the solutions regarding to the parameter t_f and we check the second-order optimality condition by a rank test to detect any conjugate points. The methods implemented in the software HAMPATH are developed more in details Ref. 15 as well as the results presented in the last section.

4.1. Homotopy and Simple Shooting in the Smooth case

In order to find the structure and a good approximation of the solutions of the contrast problem, we define a Tychonov Ref. 24 type regularization of the problem and two such regularizations are analyzed, using continuation techniques.

12 *B. BONNARD, O. COTS*

The first one which will be denoted by $L^{2-\lambda}$ -regularization transforms the Mayer problem in the Bolza form with cost:

$$c(x(t_f)) + (1 - \lambda) \int_0^{t_f} |u|^{2-\lambda} dt, \quad \lambda \in [0, 1],$$

while the second one called L^2 -regularization is:

$$c(x(t_f)) + (1 - \lambda) \int_0^{t_f} |u|^2 dt, \quad \lambda \in [0, 1].$$

If the control is saturating then its value is given by $u(t) = 2\pi \cdot \text{sgn}(H_G)$ in both cases. Otherwise, according to the maximization condition we have for the $L^{2-\lambda}$ -regularization:

$$u(x, p, \lambda) = \text{sgn}(H_G) \left(\frac{2|H_G|}{(2 - \lambda)(1 - \lambda)} \right)^{\frac{1}{1-\lambda}},$$

while for the L^2 -regularization, the control u is equal to

$$u(x, p, \lambda) = \frac{H_G}{(1 - \lambda)},$$

where $p(\cdot)$ is the adjoint vector resulting from the Pontryagin Maximum Principle. Plugging $u(x, p, \lambda)$ into the Hamiltonians describing both regularizations, we have the following Hamiltonian system:

$$\frac{dx}{dt} = \frac{\partial H}{\partial p}(x, p, \lambda), \quad \frac{dp}{dt} = -\frac{\partial H}{\partial x}(x, p, \lambda),$$

with

$$\begin{aligned} H(x, p, \lambda) &= p^0(1 - \lambda)|u(x, p, \lambda)|^{2-\lambda} + H_F(x, p) + u(x, p, \lambda) H_G(x, p) \\ &\text{or} \\ H(x, p, \lambda) &= p^0(1 - \lambda)|u(x, p, \lambda)|^2 + H_F(x, p) + u(x, p, \lambda) H_G(x, p), \end{aligned}$$

which do not depend on u anymore. We can now introduce the one-parameter family of shooting equations (or *homotopic function*), which is defined by the boundary conditions (3.8):

$$\begin{aligned} S : \Omega \subset \mathbb{R}^n \times [0, 1] &\rightarrow \mathbb{R}^n \\ \begin{pmatrix} p_0 \\ \lambda \end{pmatrix} &\mapsto S(p_0, \lambda) = \begin{pmatrix} x_1(t_f, x_0, p_0, \lambda) \\ x_2(t_f, x_0, p_0, \lambda) - p_2(t_f, x_0, p_0, \lambda) \end{pmatrix}, \end{aligned}$$

with $x_0 = x(0) = N = ((0, 1), (0, 1))$.

For each regularization, we first solve the problem in $\lambda = 0$ using a simple shooting method. We use the fortran hybrid Newton method `hybrj` (from the `minpack` library) to solve the nonlinear equation $S(p_0, 0) = 0$. The Jacobian of the restriction of S on $\lambda = 0$ is given to `hybrj` and must be invertible. Note that for $\lambda = 0$, both regularizations represent the same problem and so the solutions are similar.

If the interior $\mathring{\Omega} \subset \mathbb{R}^n \times (0, 1)$ of the domain Ω is made only of regular points of S and if the restriction of S on $\lambda = 0$ is a submersion, *i. e.*:

$$\begin{aligned} \text{rank } S'(p_0, \lambda) &= n, \quad (p_0, \lambda) \in \mathring{\Omega}, \\ \text{rank } \frac{\partial S}{\partial p_0}(p_0, \lambda)|_{\lambda=0} &= n, \quad p_0 \in \mathbb{R}^n, \end{aligned}$$

then as a consequence, the level set $\{S = 0\}$ is a one-dimensional submanifold of \mathbb{R}^{n+1} called the *path of zeros*, starting from $\lambda = 0$ to $\lambda = \lambda_f < 1$. For any $r = (p_0, \lambda) \in \Omega$, $\dim \ker S'(r) = 1$ so one can define the (tangent) vector $T(r)$ as being the unique—up to orientation—unit vector in the kernel. The orientation is chosen so that the nonvanishing determinant

$$\det \begin{bmatrix} S'(r) \\ {}^t T(r) \end{bmatrix}$$

has constant sign on each connected component of Ω . This provides a parameterization by arc length of the connected components of $\{S = 0\}$ which are computed by integrating the following differential equation (with $' = d/ds$):

$$r'(s) = T(r(s)), \quad r(0) = r_0 \in \{S = 0\},$$

with $r_0 = (p_0, 0)$ obtained by a first shooting. Unlike well-known prediction-correction methods, see Refs. 1, we merely follow the path of zeros by integrating the differential system with a high order Runge-Kutta scheme, without any correction. The solution at λ_f should be an approximation of the solution of the contrast problem, in order to identify the structure of the control and to initialize the multiple shooting method described in the next part.

The main issue here is to accurately compute the Jacobian of the homotopic function. In this case we have S of the form $S(p_0, \lambda) = b(z(t_f, x_0, p_0, \lambda)) = b(z_f)$, with $b : \mathbb{R}^{2n} \rightarrow \mathbb{R}^n$ and $z = (x, p)$. The derivative $b'(z)$ is computed in the HAMPATH code by Automatic Differentiation. The derivative $\partial_{p_0} z(t_f, x_0, p_0, \lambda)$ is the solution at t_f of the Jacobi equation:

$$\delta \dot{z} = \frac{\partial \vec{H}}{\partial z}(z(t), \lambda) \delta z, \quad (4.1)$$

with the initial value $\delta z(0) = (0, I_n) \in \mathbb{R}^{2n \times n}$. In a same way, the partial derivative with respect to the homotopic parameter λ , $\partial_\lambda z(t_f, x_0, p_0, \lambda)$, is the solution at t_f of:

$$\delta \dot{z} = \frac{\partial \vec{H}}{\partial z}(z(t), \lambda) \delta z + \frac{\partial \vec{H}}{\partial \lambda}(z(t), \lambda),$$

with the initial value $\delta z(0) = 0 \in \mathbb{R}^{2n}$. Again HAMPATH uses automatic differentiation to compute \vec{H} , $\partial_z \vec{H}$ and $\partial_\lambda \vec{H}$.

4.2. Multiple Shooting and Homotopy in the Bang-Singular case

Assume to simplify that the optimal solution of the contrast problem (*i.e.* at $\lambda = 1$) is made of a single BS-sequence and let us denote by t_1 , $0 < t_1 < t_f$, the switching time between the regular arc and the singular arc. By construction, $z(t)$, $t \in [0, t_f]$ is a concatenation of the integral curve of \vec{H} joining $z(0)$ and $z_1 = z(t_1)$, where $u(t) = 2\pi \operatorname{sgn} H_G(z(t))$ and of the integral curve of \vec{H}_s starting at z_1 , where z_1 must satisfy Eq. (3.6). Let us denote by $\tilde{z}(\cdot, z_1, t_1)$, $\tilde{z} = (\tilde{x}, \tilde{p})$ the singular extremal starting at $z(t_1) = z_1$.

We can define the function which maps (p_0, t_1, z_1) to the boundary conditions, to the Eq. (3.6) and to the matching conditions $z(t_1, x_0, p_0) = z_1$.

This leads to the multiple shooting function, see Ref. 11, 23:

$$S : \mathbb{R}^{3n+1} \rightarrow \mathbb{R}^{3n+1}$$

$$\begin{pmatrix} p_0 \\ t_1 \\ z_1 \end{pmatrix} \mapsto \begin{pmatrix} H_G(z_1) \\ \dot{H}_G(z_1) \\ z(t_1, x_0, p_0) - z_1 \\ \tilde{x}_1(t_f, z_1, t_1) \\ \langle \tilde{x}_2(t_f) | \tilde{p}_2(t_f) \rangle - |\tilde{x}_2(t_f)|^2 \end{pmatrix},$$

which is of the form $S(p_0, t_1, z_1) = b(t_1, z_1, z(t_1, x_0, p_0), \tilde{z}(t_f, z_1, t_1))$. We again use a Newton type algorithm to solve the nonlinear equation $S(p_0, t_1, z_1) = 0$. We need the Jacobian of S , which is quite different from the smooth case. The partial derivatives of b are more intricate than before and are computed again using Automatic Differentiation. As in the smooth case, $\partial_{p_0} z(t_1, x_0, p_0)$ and $\partial_{z_1} \tilde{z}(t_f, z_1, t_1)$ are solutions of the Jacobi equations (4.1), respectively with the Hamiltonian vector fields \vec{H} and \vec{H}_s . Then $\partial_{t_1} z(t_1, x_0, p_0)$ is simply equal to $\vec{H}(z(t_1, x_0, p_0))$ and finally $\partial_{t_1} \tilde{z}(t_f, z_1, t_1)$ is the solution at t_f of:

$$\delta \dot{z} = \frac{\partial \vec{H}_s}{\partial z}(z(t, z_1, t_1)) \delta z,$$

with the initial value $\delta z(t_1) = -\vec{H}_s(z_1) \in \mathbb{R}^{2n}$.

Up to this point, the final time t_f was fixed. One important problem is to study the behavior of the solutions regarding to the parameter t_f . This is done by changing the parameterization of our problem so that our new time \bar{t} is in $[0, 1]$. One can introduce $t_f \cdot \bar{t} = t$ which leads to consider t_f as an homotopic parameter. The Hamiltonians are multiplied by t_f and we have a one-parameter family of multiple shooting equations of the form $S(p_0, \bar{t}_1, z_1, t_f)$, $\bar{t}_1 = t_1/t_f$, which can be solved by continuation techniques presented in 4.1.

4.3. Second-Order Sufficient Optimality conditions

The second order sufficient conditions of optimality are related to the concept of conjugate point. According to definition 3.2, a time $t_c \in (t_1, t_f]$, is a conjugate time if there exists a Jacobi field $\delta z(\cdot) = (\delta q(\cdot), \delta p(\cdot))$, semi-vertical at t_1 and t_c . Hence,

a Jacobi field along a singular extremal, semi-vertical at t_1 must satisfy these three conditions:

(i)

$$\begin{aligned} dH_G(z(t_1)).\delta z(t_1) &= -\delta z_1.p_{y_1} + \delta y_1.p_{z_1} - \delta z_2.p_{y_2} + \delta y_2.p_{z_2} \\ &\quad - z_1.\delta p_{y_1} + y_1.\delta p_{z_1} - z_2.\delta p_{y_2} + y_2.\delta p_{z_2} \\ &= 0 \end{aligned}$$

$$\begin{aligned} d\{H_F, H_G\}(z(t_1)).\delta z(t_1) &= -\delta_1.\delta z_1.p_{y_1} - \delta_1.\delta y_1.p_{z_1} - \delta_2.\delta z_2.p_{y_2} - \delta_2.\delta y_2.p_{z_2} \\ &\quad + (\gamma_1 - \delta_1.z_1).\delta p_{y_1} - \delta_1.y_1.\delta p_{z_1} + (\gamma_2 - \delta_2.z_2).\delta p_{y_2} - \delta_2.y_2.\delta p_{z_2} \\ &= 0; \end{aligned}$$

(ii) $\delta x(t_1) \in \text{vect}\{G(x(t_1))\}$;

(iii) $p(t_1)\delta p(t_1) = 0$ (linearization of $|p(t_1)| = 1$ which breaks the homogeneity in (p^0, p)), which means $[0_{1 \times n} \quad {}^t p(t_1)].\delta z(t_1) = 0$.

The second condition can be rewritten as follow. Since $G(x(t_1))$ is non-zero, there exists $A \in L(\mathbb{R}^n, \mathbb{R}^{n-1})$ of maximal rank $(n-1)$ such as $\text{vect}\{G(x(t_1))\} = \ker A$ ($A \in \mathbb{R}^{(n-1) \times n}$ is formed by the vectors of a basis of $\{G(x(t_1))\}^\perp$); let us denote $B = [A \quad 0_{(n-1) \times n}] \in L(\mathbb{R}^{2n}, \mathbb{R}^{n-1})$, then $\delta x(t_1) \in \text{vect}\{G(x(t_1))\}$ is equivalent to $B.\delta z(0) = 0$.

Finally if we denote by C the matrix:

$$C = \begin{pmatrix} 0_{1 \times n} \quad {}^t p(t_1) \\ dH_G(z(t_1)) \\ d\{H_F, H_G\}(z(t_1)) \\ B \end{pmatrix} \in \mathbb{R}^{(n+2) \times 2n}$$

then all the three conditions are $C.\delta z(0) = 0$. Under our assumptions C is of maximal rank, then it has a kernel of dimension $2n - (n+2) = n-2$, which defines two Jacobi fields $\delta z_i(t) = (\delta x_i(t), \delta p_i(t))$, $i = 1, 2$, if $n = 4$.

A time $t_c > t_1$ is a conjugate time if we have the following rank condition:

$$\text{rank}\{\delta x_1(z(t_c)), \delta x_2(z(t_c)), G(x(t_c))\} < 3.$$

Using assumption (A3), this is equivalent to:

$$\det\{\delta x_1(z(t_c)), \delta x_2(z(t_c)), G(x(t_c)), F(x(t_c))\} = 0. \quad (4.2)$$

5. Numerical results in the case of blood samples

5.1. Comparison of the two homotopies

When λ is equal to zero, the two homotopies coincide; remarkably, both are initialized by using a trivial guess for the unknown initial adjoint state, $p_0 = 5e - 2 \times (1, 1, 1, 1)$. The solution Fig. 3 for $\lambda = 0$ is obtained with an error of 10^{-14} on the norm of the shooting function. The path of zeros is then followed up

16 *B. BONNARD, O. COTS*

to some $\lambda_f < 1$, revealing the BS structure of the control (see Figs. 4–5). The evolution of the contrast along the path is given Fig. 6. The path itself is represented in Fig. 7. All results of this subsection and the following are obtained for t_f fixed to 1.1 times the minimum time of the single spin system with the same boundary conditions as spin no. 1.

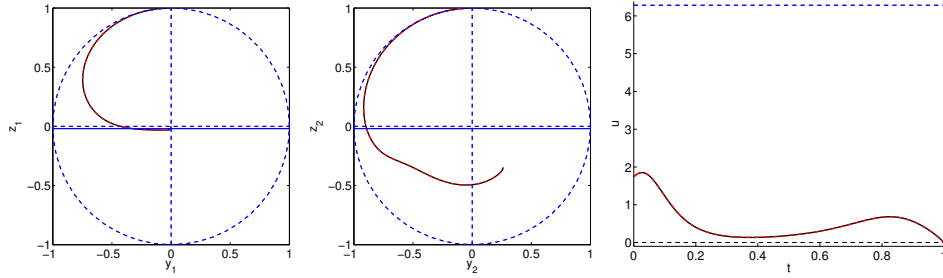


Fig. 3. Solution for $\lambda = 0$ ($t_f = 1.1 \times \min t_f$). Trajectories for spin 1 and 2 in the (y, z) -plane are portrayed in the first two subgraphs. In each subgraph, the solution associated with the L^2 -homotopy (*resp.* $L^{2-\lambda}$) is the dashed red line (*resp.* black line). The corresponding controls are drawn in the rightmost subgraph. For $\lambda = 0$, the two homotopies define the same problem, so trajectories and controls are identical for both.

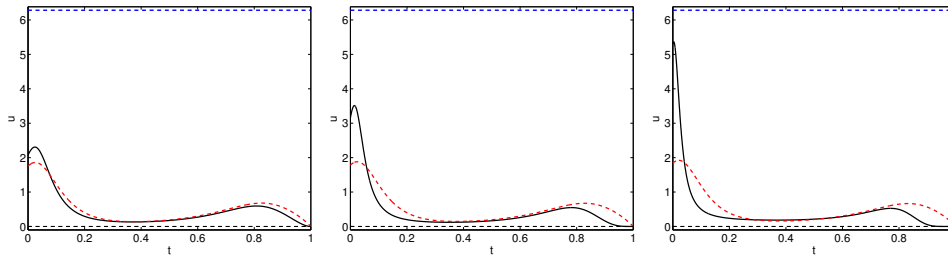


Fig. 4. Controls for $\lambda = 0.5, 0.8, 0.9$ ($t_f = 1.1 \times \min t_f$). In each subgraph, the solution associated with the L^2 -homotopy (*resp.* $L^{2-\lambda}$) is the dashed red line (*resp.* black line). As $\lambda \rightarrow 1$, the BS structure is numerically revealed for by the $L^{2-\lambda}$ -homotopy. It is less visible for the L^2 -homotopy.

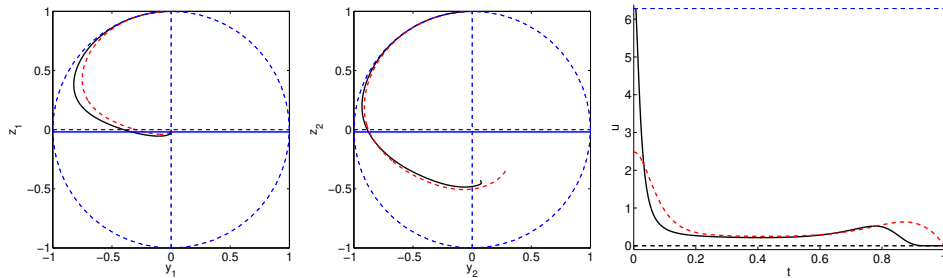


Fig. 5. Solutions for $\lambda_f < 1$ ($t_f = 1.1 \times \min t_f$). Trajectories and controls as in Fig.3. In the L^2 case, $\lambda_f = 0.99$; in the $L^{2-\lambda}$ case, $\lambda_f = 0.92$.

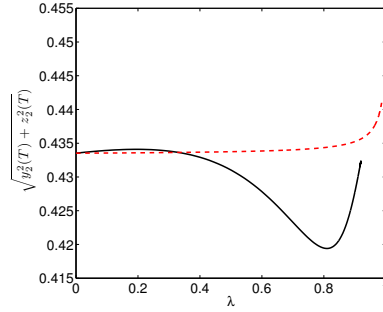


Fig. 6. Contrast along the path ($t_f = 1.1 \times \min t_f$). The contrast at the final state is evaluated for each homotopy as λ ranges $[0, \lambda_f]$. The value associated with the L^2 -homotopy (*resp.* $L^{2-\lambda}$) is the dashed red line (*resp.* black line). Although it is not clear whether the contrasts associated with each one of the two homotopies converge to the same value as $\lambda \rightarrow 1$ (see also Fig. 7), the corresponding solutions allow to initialize the resolution for $\lambda = 1$ and lead then to the same solution of the contrast problem (see §5.2).

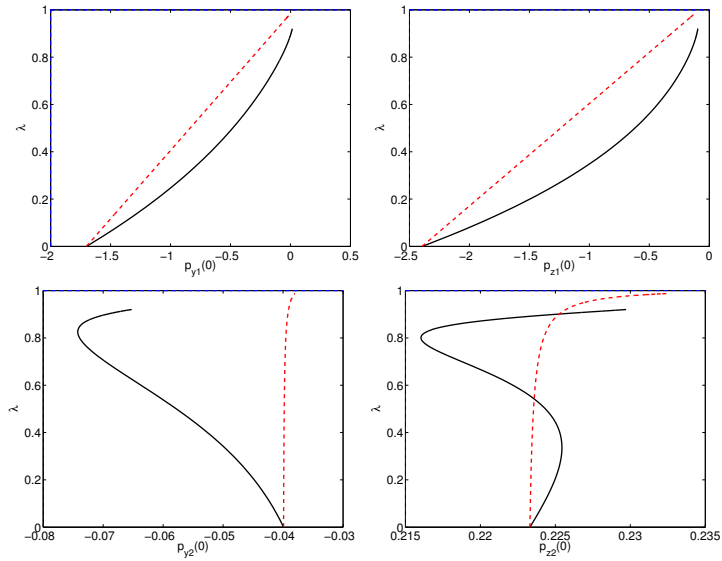


Fig. 7. Path of zeros ($t_f = 1.1 \times \min t_f$). For each component of the unknown initial costate, the path associated with the L^2 -homotopy (*resp.* $L^{2-\lambda}$) is the dashed red line (*resp.* black line). The main differences when $\lambda \rightarrow 1$ are observed on p_{y_2} and p_{z_2} .

5.2. Resolution of the contrast problem

The value obtained at $\lambda_f < 1$ for each homotopy is a candidate as initial guess for the multiple shooting when one assumes the BS structure. One also needs to initialize the unknown junction time t_1 between the bang and singular arc. A whole range of initializations is considered as illustrated by Fig. 8. These various initializations lead to three different solutions (see Figs. 9–11), no matter the homotopy chosen to initialize p_0 . The influence of the initial guess for t_1 on the result is discussed Fig. 12. The error on the norm of the shooting function is at most equal to 10^{-10} for any solution.

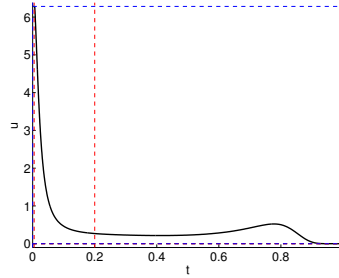


Fig. 8. Discretization of the time interval to define initial guesses for t_1 ($t_f = 1.1 \times \min t_f$). The initializations of the unknown junction time t_1 from the bang to the singular arc are chosen in the subinterval $[0.005, 0.2]$. This heuristic choice is related to the structure of the control for λ_f (here, the control obtained for the $L^{2-\lambda}$ homotopy is portrayed). The corresponding initialization for $z_1 = z(t_1)$ is obtained by merely integrating the dynamics for λ_f .

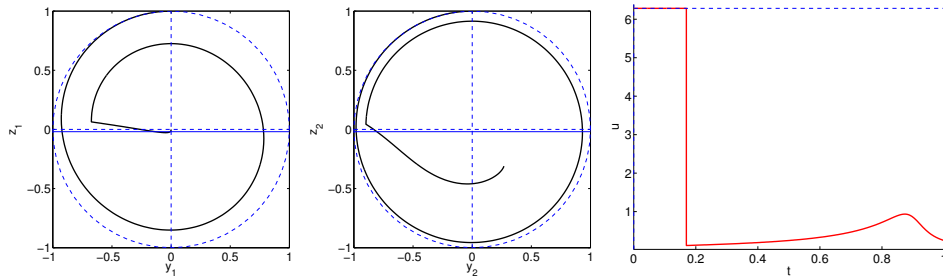


Fig. 9. Contrast problem, solution no. 1 with contrast 0.411 ($t_f = 1.1 \times \min t_f$). Trajectories for spin 1 and 2 in the (y, z) -plane are portrayed in the first two subgraphs. The corresponding BS control is drawn in the rightmost subgraph.

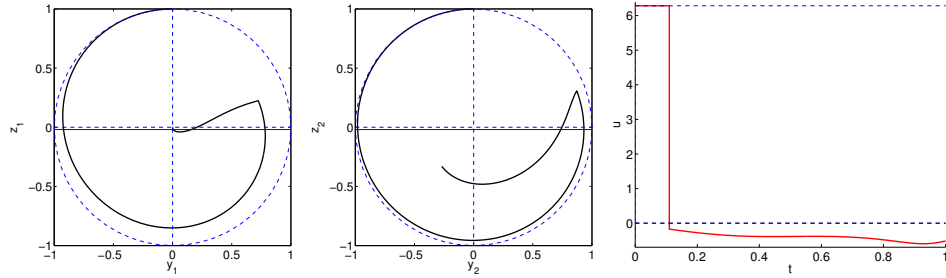


Fig. 10. Contrast problem, solution no. 2 with contrast 0.426 ($t_f = 1.1 \times \min t_f$). Trajectories for spin 1 and 2 in the (y, z) -plane are portrayed in the first two subgraphs. The corresponding BS control is drawn in the rightmost subgraph.

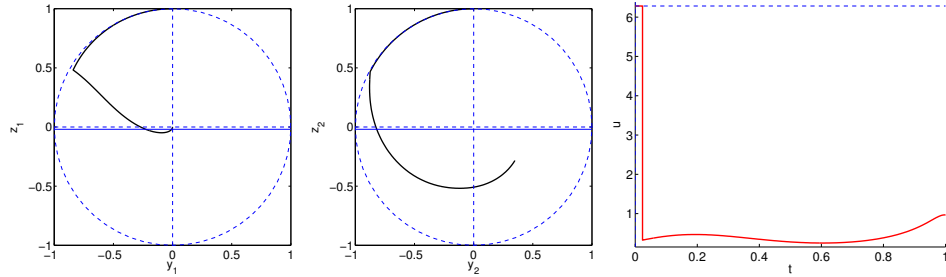


Fig. 11. Contrast problem, solution no. 3 with contrast 0.449 ($t_f = 1.1 \times \min t_f$). Trajectories for spin 1 and 2 in the (y, z) -plane are portrayed in the first two subgraphs. The corresponding BS control is drawn in the rightmost subgraph.

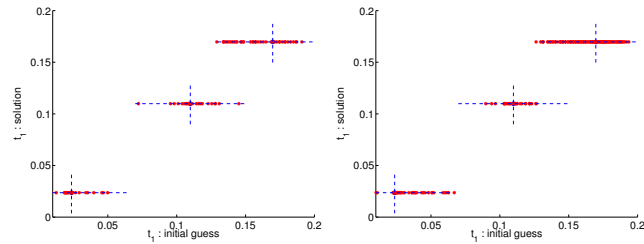


Fig. 12. Influence of the initial guess for t_1 on the contrast solution ($t_f = 1.1 \times \min t_f$). For each value of the initial guess on the x -axis, the solution for t_1 is plotted when convergence is obtained. The initial costate p_0 is either initialized for the solution for λ_f of the L^2 - λ -homotopy (LHS), or the L^2 -homotopy (RHS). As is clear from Figs. 9, 10, 11, the smaller t_1 , the better the contrast: Solutions no. 1, 2, 3 are obtained for $t_1 = 0.17, 0.11$ and 0.024 , respectively.

20 *B. BONNARD, O. COTS*

5.3. Influence of the final time

Given that the initial point is a stationary point, the contrast is an increasing function of t_f acting as a parameter. Indeed, applying a zero control at $t = 0$ leaves the system in its initial state so there is an inclusion of admissible controls between problems when the final time is increased (and the bigger the set of controls, the larger the maximum contrast). Whereas the branches no. 1 and 2 generated by the solutions for $t_f = 1.1 \times \min t_f$ provide local but not global minima, it is not possible to decide whether branch 3 corresponds to global minimizers. Accordingly, even branch 3 only provides a lower bound for the value function (see Fig. 13). Having increasing bounded (by one, which is the maximum possible contrast given the final condition on spin no. 1) functions, it is natural to expect asymptotes on each branch. The numerical simulations indicate that the limit value on the three branches is almost reached past $t_f \simeq 1.3 \times \min t_f$. For large enough final times, solutions 1 and 2 coincide for symmetry reasons (see Figs. 18–19) while the bang arc tends to disappear on solution 3 (see Fig. 20). Conversely, a BSBS structure seems to be revealed on the third branch as $t_f \rightarrow \min t_f$ (see Figs. 14–17). The corresponding three paths of zeros are represented on Fig. 21. Second order condition checks are also performed along branch 3 as illustrated Fig. 22, ensuring local optimality by virtue of Theorem 3.2.

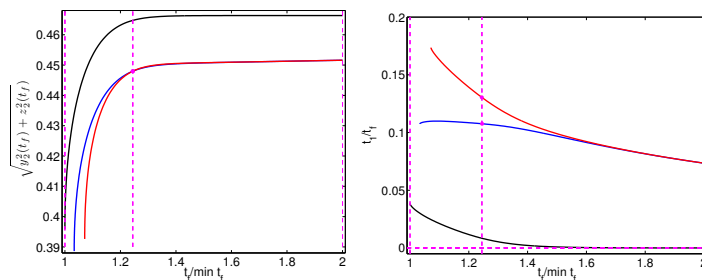


Fig. 13. Evolution of the contrast with respect to the final time. On the left, the contrast increases as t_f is varied up to two times the minimum time of the one spin system for solution no. 1 (red line), 2 (blue) and 3 (black line). For $t_f > 1.3 \times \min t_f$, the asymptotic value on each branch of solutions is almost reached (up to 99.5% of the maximal contrast). Moreover, although the costs associated with branches no. 1 and 2 are indiscernible past such final times, the strategies still differ as illustrated on the rightmost subgraph (see also Figs. 18–19). At $1.244 \times \min t_f$, the costs cross and then tend to the same value for larger t_f . On the right, the evolution of the junction time t_1 is portrayed for each branch of solution. On one hand, for $t_f \simeq 1.244 \times \min t_f$, the values of t_1 still differ between branches no. 1 and 2. This difference tends to vanish for larger t_f . On the other hand, $t_1 \rightarrow 0$ on branch 3 as t_f grows, entailing that the bang arc tends to disappear.

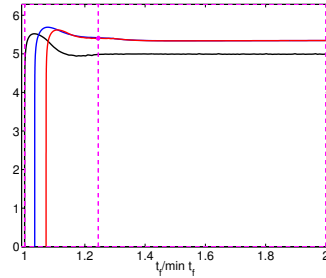


Fig. 14. The minimal distance between the singular control along the trajectory and the upper bound 2π with respect to the normalized final time, for solution no. 1 (red line), 2 (blue) and 3 (black line). The distance tends to 0 when t_f decreases. When it is 0, the singular control is saturating and another bang arc is required. For solutions no. 1 and 2, the distance is 0 for t_f greater than $\min t_f$. On the other hand, for solution no. 3 the distance tends to 0 as t_f tends to $\min t_f$ and the BSBS-sequence of the single spin case is revealed (see Fig. 18).

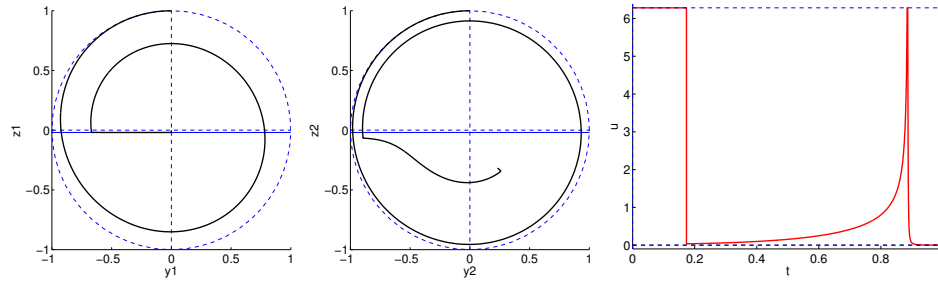


Fig. 15. Solution no. 1 ($t_f = 1.07 \times \min t_f$). Trajectories for spin 1 and 2 in the (y, z) -plane are portrayed in the first two subgraphs for solution no. 1. The corresponding control is drawn in the rightmost subgraph. As $t_f \rightarrow \min t_f$, the BSBS structure is revealed as the singular subarc comes closer to saturating the constraint $|u| \leq 2\pi$.

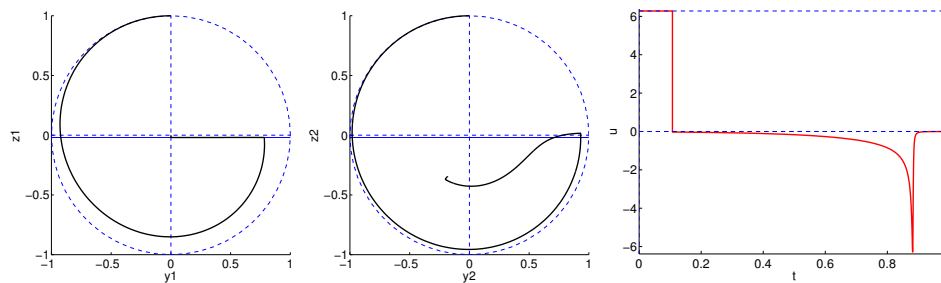


Fig. 16. Solution no. 2 ($t_f = 1.03 \times \min t_f$). Idem as Fig. 15.

22 *B. BONNARD, O. COTS*

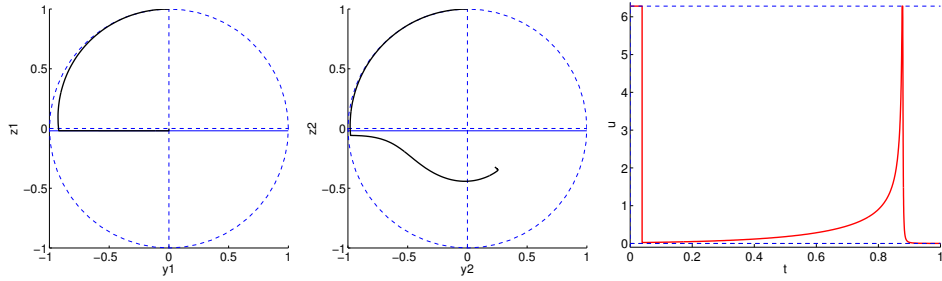


Fig. 17. Solution no. 3 ($t_f = 1.000004 \times \min t_f$). Idem as Fig. 15.

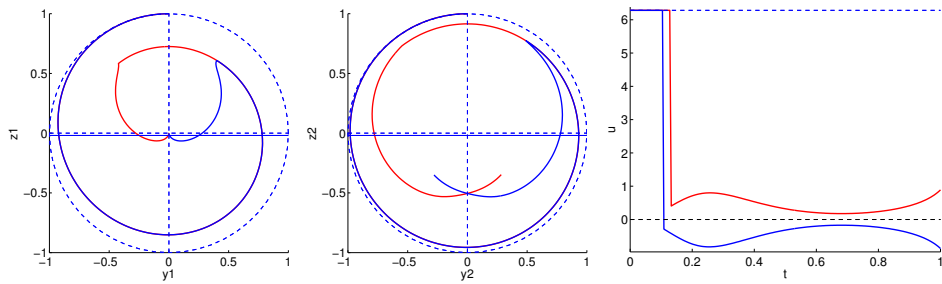


Fig. 18. Solutions no. 1 and 2 ($t_f = 1.244 \times \min t_f$). Trajectories for spin 1 and 2 in the (y, z) -plane are portrayed in the first two subgraphs for solution no. 1 (red line) and 2 (blue line). The corresponding controls are drawn in the rightmost subgraph. Despite the small difference between the durations of the bang arcs, the difference in strategies is clear and associated with a symmetry of the system.

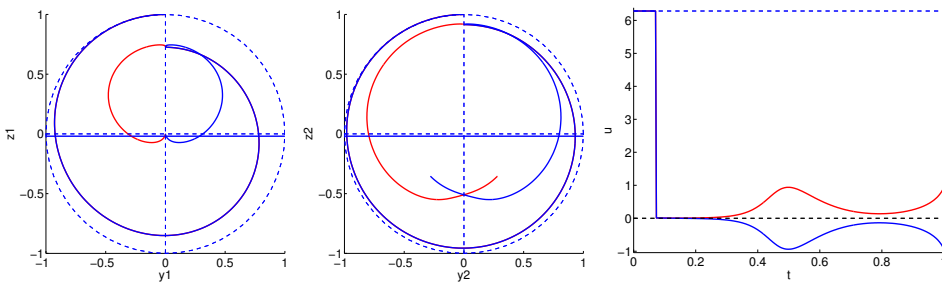


Fig. 19. Solutions no. 1 and 2 ($t_f = 2 \times \min t_f$). Trajectories for spin 1 and 2 in the (y, z) -plane are portrayed in the first two subgraphs for solution no. 1 (red line) and 2 (blue line). The durations of the bang arcs are equal and the singular controls are opposite. The singular trajectories are symmetrical with respect to the z -axis.

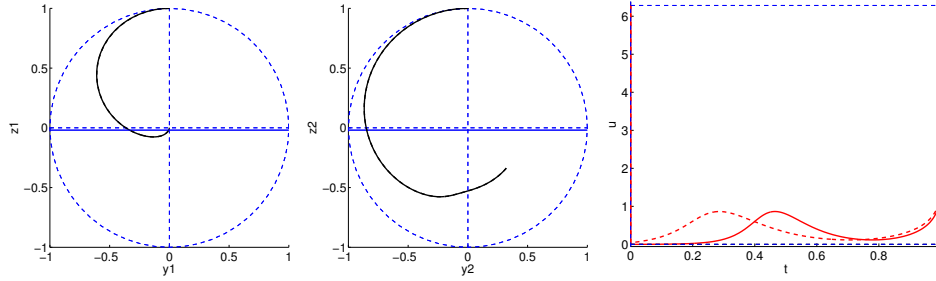


Fig. 20. Solution no. 3: $t_f = 1.5 \times \min t_f$ (dashed line) and $t_f = 2 \times \min t_f$ (filled line). Trajectories for spin 1 and 2 in the (y, z) -plane are portrayed in the first two subgraphs for solution no. 3. The corresponding control is drawn in the rightmost subgraph. The trajectories are superimposed but not the controls. The bang arc tends to disappear and the singular control makes the particles resting at the initial point close to the north pole during more time.

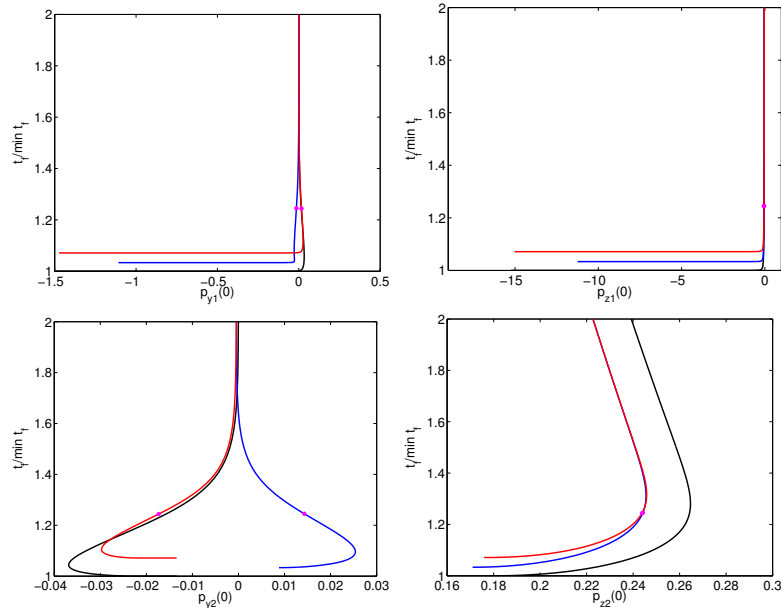


Fig. 21. Path of zeros for the continuation on the final time. For each component of the unknown initial costate, branches associated with solution no. 1 (red line), 2 (blue line) and 3 (black line) are portrayed. The convergence of branches 1 and 2 is observed as t_f increases, in coherence with the convergence on t_1 detailed Fig. 13. The only persisting difference with branch no. 3 comes from the p_{y_2} and p_{z_2} components of the initial adjoint state.

24 B. BONNARD, O. COTS

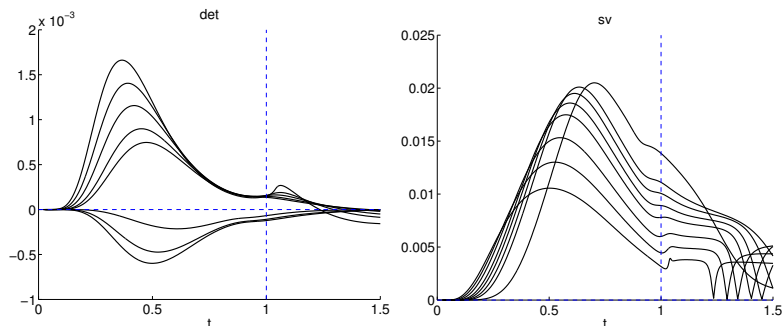


Fig. 22. Second order sufficient condition check on branch no. 3. The rank condition resulting from Theorem 3.2 is evaluated for several values of the ratio $t_f / \min t_f$ in $(1, 1.15]$ for the branch of solutions no. 3 that provide the best contrast. On the left, for each t_f , the curve is reparameterized so that the final time corresponds to the abscissa 1; the determinant associated with the rank condition is plotted, so there is a conjugate time whenever it vanishes. One observes that conjugate times on each extremal are located after the (normalized to 1) final time, ensuring local optimality of the trajectory. On the right, the smallest singular value of the matrix whose rank we test is plotted, extracting only the relevant information to detect the rank drops (whereas the previous determinant is equal to the product of all singular values).

6. Conclusion

In this article, we have presented the geometric optimal methods combined with numerical adapted schemes implemented in the HAMPATH code which lead to calculate the optimal control in the NMR-imaging techniques for the blood sample case. Restricting to a single control field, a contrast of 0.466 was obtained for a BS-sequence and a transfer time of $1.5 \times \min t_f$, where $\min t_f$ is the minimum time to drive the first spin to zero magnetization. This gives a local optimum with respect to all solutions contained in a tubular neighborhood of the singular arc.

The same techniques can be applied to other cases, e.g. cerebrospinal fluid/water samples. In the analysis presented in this article, the continuation method is applied to regularize the Mayer problem to Bolza and a second continuation concerns the transfer time. Similarly a continuation method can be made on the relaxation parameters to study the different cases.

As a conclusion, to illustrate the extension of the approach we present a numerical result. A BSBS-sequence, see Fig. 23, is computed thanks to the regularization process in the cerebrospinal fluid/water case and a continuation is made towards the blood case which leads to a change of the structure in BSBSBS, see Fig. 24, and provides a contrast of 0.484 for the 3BS-sequence for a transfer time of $1.5 \times \min t_f$ vs. 0.466 for a BS-sequence of the blood case analyzed in this article.

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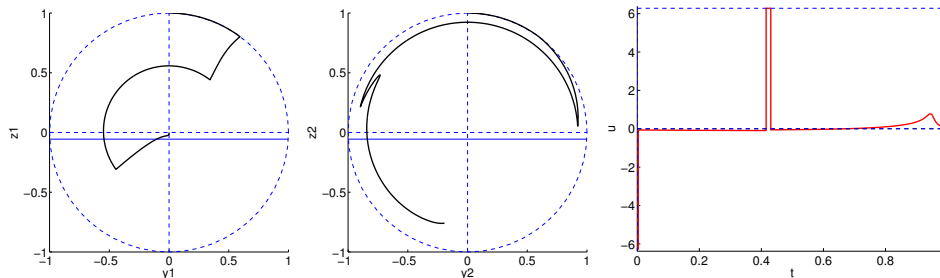


Fig. 23. Contrast problem in the fluid/water case, solution BSBS with contrast 0.783 for a transfer time $t_f = 1.5 \times \min t_f$. Trajectories for spin 1 (*i.e.* cerebrospinal fluid with $T_1 = 2000\text{ms}$ and $T_2 = 200\text{ms}$) and spin 2 (*i.e.* water with $T_1 = 2500\text{ms}$ and $T_2 = 2500\text{ms}$) in the (y, z) -plane are portrayed in the first two subgraphs. The corresponding control is drawn in the rightmost subgraph.

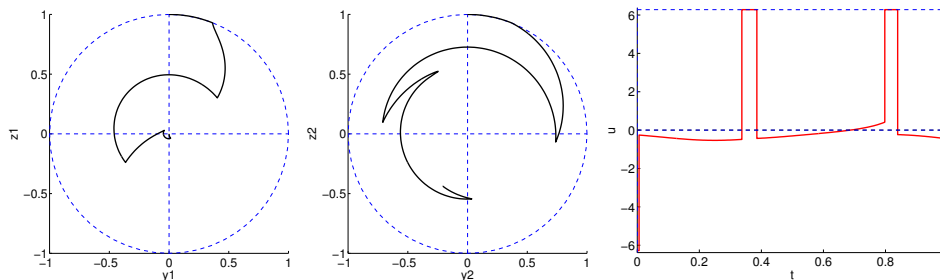


Fig. 24. Contrast problem in the blood case, solution BSBSBS with contrast 0.484 for a transfer time $t_f = 1.5 \times \min t_f$, from continuation from the fluid/water case. Trajectories for spin 1 (*i.e.* deoxygenated blood with $T_1 = 1350\text{ms}$ and $T_2 = 50\text{ms}$) and spin 2 (*i.e.* oxygenated blood $T_1 = 1350\text{ms}$ and $T_2 = 200\text{ms}$) in the (y, z) -plane are portrayed in the first two subgraphs. The corresponding control is drawn in the rightmost subgraph.

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26 *B. BONNARD, O. COTS*

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