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WORKING NOTES ON THE TIME MINIMAL SATURATION OF
A PAIR OF SPINS AND APPLICATION IN MAGNETIC
RESONANCE IMAGING

B. BONNARD, O. COTS, J. ROUOT, AND T. VERRON

ABSTRACT. In this article, we analyse the time minimal control for the satu-
ration of a pair of spins of the same species but with inhomogeneities of the
applied RF-magnetic field, in relation with the contrast problem in Magnetic
Resonance Imaging. We make a complete analysis based on geometric control
to classify the optimal syntheses in the single spin case to pave the road to
analyze the case of two spins. The Bocop software is used to determine local
minimizers for physical test cases and Linear Matrix Inequalities approach is
applied to estimate the global optimal value and validate the previous comput-
tations. This is complemented by numerical computations combining shooting
and continuation methods implemented in the HamPath software to analyze the
structure of the time minimal solution with respect to the set of parameters
species.

1. Introduction

In Magnetic Resonance Imaging (MRI), a challenging problem is to maximize
the contrast between two observed species, e.g. healthy tissues from tumors by
saturating one of the species. Optimal control techniques were introduced in this
domain in the eighties [17] and were developed very recently under the impulse
of S. Glaser using advanced analytical and numerical techniques. This gave rise
to a series of articles which analyze the contrast problem, starting from the ideal
contrast problem where only a pair of spins is considered, to the optimal control
of an ensemble of pairs of spins, taking into account the so-called $B_0$ and $B_1$
inhomogeneities [22, 23, 11, 8] to compute robust control. A first major contribution in
this area was for some physical cases, the explicit computation of the time minimal
control of saturating a single spin, that is steering the state representing the spin
to zero and showing that the standard inversion sequence applied in practice is not
optimal in many physical cases, e.g. the blood case [22]. Additional pulses have to
be used corresponding to the so-called singular control [5]. This computation relies
on an intense research activities at the end of the eighties to the computation of
the time-minimal closed loop solution for $C^\omega$-planar single input control system in
a neighborhood of a given point, taking into account the Lie algebraic structure of
the systems of this point [12, 35, 33, 34].

Thanks to the bilinear structure of the model, this local analysis can be global-
ized for a single spin and a first contribution of our article is to make a complete
classification of the time minimal syntheses taking into account the time relaxation
parameters of the species and the maximal amplitude of the applied RF-field (con-
trol). Our aim is to extend this analysis to the case of two spins in relation to the
problem of the so-called $B_1$-inhomogeneity, that is of the applied RF-field. The
analysis is very complex and we present a combination of algebraic and geometric
methods based on [6] and adapted numerical schemes to analyze the problem im-
plemented in specific software: Bocop [3], HamPath [15], GloptiPoly [19]. A similar
analysis was presented for the contrast problem [11] but for this simplified version
it is complete in many directions, thus providing a testbed of the methods to deal
with the time minimal control of a single input control system where the state space
is of codimension 4, depending upon 2-parameters and with many local optima.

This article is organized as follows. In section 2, we present the mathematical
model, that is the Bloch equations [26] and we discuss the underlying optimal con-
trol problem in MRI, that is the contrast problem with $B_0$ and $B_1$ inhomogeneities
[23], to introduce the time minimal saturation of a pair of spins, that we ana-
lyze in this article. The seminal result in optimal control theory is the Maximum
Principle [31] which is recalled to select extremal curves candidates as minimizers.
The extremal controls split into bang controls and the so-called singular controls
whose role in the time minimal problem is recalled [7]. In section 3, the time
minimal saturation of a pair of spins of the same species with $B_1$-inhomogeneities
is investigated. A preliminary study concerning the case of a simple spin is pre-
sented in details using geometric control theory techniques. This led to a complete
classification of the optimal syntheses to steer any state to the origin. Thanks to
the symmetry of revolution it is reduced to a time minimal control problem for a
single-input 2D-system. All the fine results of the geometric theory [13, 35] are used
to provide a complete classification depending upon the physical parameters and
completing [22]. The next step in section 4 is to extend this analysis to a pair of
spins, numerical methods are used to complete the analysis. First of all, the direct
methods implemented in the Bocop code are applied to analyze for physical cases:
Desoxygenated and Oxygenated case, Cerebrospinal fluid and Water case. Global
optimality is analyzed using LMI methods [24]. In the final part, the problem is an-
alyzed using multiple shooting methods implemented in the HamPath software [15]
and completed by numerical continuations (available in the software) to compute
the optimal solutions for a continuous set of physical relaxation parameters. This
numerical investigation, based on homotopy reveals the existence path of zeros that
has to be compared to determine the global optimal. In section 5, we discuss the
theoretical complexity of the procedure. The crucial point is to analyze the singular
trajectories associated with a 4D-Hamiltonian flow, with constraints and many sin-
gularities. They are computed using symbolic computations, extending techniques
from [6]. They led to identify a simplified case corresponding to the so-called water
case, important in practice and which is a generalization of the standard inversion
sequence for a single spin. This case is important to analyze the general case, using
homotopy methods.
Table 1. Matter name with relaxation times in seconds, ratio $T_2/T_1$ and value $\theta = \atan(\gamma/\Gamma)$.

<table>
<thead>
<tr>
<th>Name</th>
<th>$T_1$</th>
<th>$T_2$</th>
<th>$\frac{T_2}{T_1}$</th>
<th>$\theta$</th>
</tr>
</thead>
<tbody>
<tr>
<td>Water</td>
<td>2.5</td>
<td>2.5</td>
<td>1.0</td>
<td>0.7854</td>
</tr>
<tr>
<td>Fat</td>
<td>0.2</td>
<td>0.1</td>
<td>0.5</td>
<td>0.4636</td>
</tr>
<tr>
<td>Cerebrospinal Fluid</td>
<td>2.0</td>
<td>0.3</td>
<td>0.15</td>
<td>0.1489</td>
</tr>
<tr>
<td>Oxygenated blood</td>
<td>1.35</td>
<td>0.2</td>
<td>0.1481</td>
<td>0.1471</td>
</tr>
<tr>
<td>White cerebral matter</td>
<td>0.78</td>
<td>0.09</td>
<td>0.1154</td>
<td>0.1148</td>
</tr>
<tr>
<td>Gray cerebral matter</td>
<td>0.92</td>
<td>0.1</td>
<td>0.1087</td>
<td>0.1083</td>
</tr>
<tr>
<td>Brain</td>
<td>1.062</td>
<td>0.052</td>
<td>0.0490</td>
<td>0.0489</td>
</tr>
<tr>
<td>De-oxygenated blood</td>
<td>1.35</td>
<td>0.05</td>
<td>0.0370</td>
<td>0.0370</td>
</tr>
<tr>
<td>Parietal muscle</td>
<td>1.2</td>
<td>0.029</td>
<td>0.0242</td>
<td>0.0242</td>
</tr>
</tbody>
</table>

2. Theoretical concepts and results

2.1. The model. We consider an ensemble of spin-1/2 particles, excited by a radio-frequency (RF) field which is ideally assumed homogeneous, each spin of this ensemble being described by the magnetization vector $M = (M_x, M_y, M_z)$ whose dynamics is governed in a specific rotating frame by the Bloch equation [26]:

$$
\dot{M}_x(t) = -\frac{M_x}{T_2} + \omega_y M_z,
$$

$$
\dot{M}_y(t) = -\frac{M_y}{T_2} - \omega_x M_z,
$$

$$
\dot{M}_z(t) = \frac{M_0 - M_z}{T_1} - \omega_y M_x + \omega_x M_y,
$$

where $T_1$, $T_2$ are respectively the longitudinal, transversal relaxation constants, $M_0$ is the thermal equilibrium and $\omega = (\omega_x, \omega_y)$ is the control corresponding to the applied RF-magnetic field, with $\omega_{\text{max}}$ the maximal amplitude of the control, i.e. $\omega_x^2 + \omega_y^2 \leq \omega_{\text{max}}^2$. Table 1 gives a list of longitudinal and transversal relaxation constants for the main practical cases.

Up to a renormalization of $M$ introducing $(x, y, z) = (M_x, M_y, M_z)/M_0$ and a time reparameterization, the dynamics take the form:

$$
\dot{x}(t) = -\Gamma x + u_y z,
$$

$$
\dot{y}(t) = -\Gamma y - u_x z,
$$

$$
\dot{z}(t) = \gamma(1 - z) - u_y x + u_x y.
$$

In the relevant physical cases, one has $0 \leq \gamma \leq 2\Gamma$ and the Bloch ball: $x^2 + y^2 + z^2 \leq 1$, is invariant for the dynamics. Thanks to the reparameterization, one can assume...
that the control is bounded by $u_2^2 + u_2^2 \leq 1$. In this case, the relations between the parameters $(\gamma, \Gamma)$ and $(T_1, T_2, \omega_{\text{max}})$ are the following:

$$\gamma = \frac{1}{T_1 \omega_{\text{max}}}, \quad \Gamma = \frac{1}{T_2 \omega_{\text{max}}}.$$  

(2.1)

The system admits a symmetry of revolution around the z-axis, which allows us to set $u_y = 0$ and to restrict each spin system to a single-input control system in the trace of the Bloch ball on the plane $(y, z)$. The system then takes the form:

$$\dot{y}(t) = -\Gamma y - uz,$$

$$\dot{z}(t) = \gamma(1 - z) + uy,$$

(2.2)

with $u = u_z$, $|u| \leq 1$.

The problem of saturation associated to the contrast problem by saturation in MRI is to steer from the north pole $N = (0, 1)$ to the origin $O = (0, 0)$ one of the two species to be distinguished. In the contrast problem with RF-inhomogeneities, which is due to the spatial position of the species in the image, one has to consider an ensemble of pair of spins, such as for each system, the dynamics is perturbed. This perturbation is modeled as a rescaling of the maximal amplitude perceived by the spin. Restricting again to the sub-problem of saturation of one species, and considering only an ensemble of two pairs of spins, this leads to consider the case of a couple of systems (2.2) with the same parameters $(\gamma, \Gamma)$ but with a distortion in the maximal amplitude, that is:

$$\begin{cases}
\dot{y}_1(t) = -\Gamma y_1 - uz_1, \\
\dot{y}_2(t) = -\Gamma y_2 - (1 - \varepsilon) u z_2,
\end{cases}$$

$$\begin{cases}
\dot{z}_1(t) = \gamma(1 - z_1) + uy_1, \\
\dot{z}_2(t) = \gamma(1 - z_2) + (1 - \varepsilon) u y_2,
\end{cases}$$

where $|u| \leq 1$, with $q_1 = (y_1, z_1)$, $q_2 = (y_2, z_2)$ denote the coordinates of spin 1 and spin 2 and $(1 - \varepsilon)$, $\varepsilon > 0$ small, is the rescaling factor of the control maximal amplitude. Hence, the saturation problem of a pair of spins consists into a simultaneous steering of the couple from $q_1(0) = q_2(0) = N$ to the center $q_1(t_f) = q_2(t_f) = O$, where $t_f$ is the transfer time. The optimal control that we shall analyze is the time minimal saturation, i.e. we aim to minimize the transfer time $t_f$.

### 2.2. Maximum principle and singular extremals.

#### 2.2.1. Preliminaries. In this section, we consider a single-input control system:

$$\frac{dq}{dt} = F + uG,$$

where $F$, $G$ are $\mathcal{C}^\omega$ vector fields defined on an open subset $V \subset \mathbb{R}^n$ and the control $u$ is a bounded measurable mapping defined on $[0, T(u)]$ and valued in $|u| \leq 1$. For fixed $q_0$ and $T > 0$, the extremity mapping is the map $E: u \in L^\infty([0, T]) \mapsto E(u) = q(T, q_0, u)$, where $q(\cdot, q_0, u)$ is the solution of the system with $q(0, q_0, u) = q_0$. A control $u \in L^\infty([0, T])$ is called singular if the extremity mapping is not of full rank and the corresponding trajectory is called singular on $[0, T]$. We have the following relations with the time minimal control problem $[31]$.

**Proposition 2.1.** Consider the time minimal control problem for the single-input control system: $\frac{dq}{dt} = F + uG$, $|u| \leq 1$. If $q(\cdot)$, with corresponding trajectory $q(\cdot)$, is solution, then there exists $p(\cdot)$, $t \mapsto p(t) \in \mathbb{R}^n \setminus \{0_{\mathbb{R}^n}\}$, such that the following
equations are satisfied for the triplet $(q(\cdot), p(\cdot), u(\cdot))$:

\begin{align}
\dot{q} &= \frac{\partial H}{\partial p}, \quad \dot{p} = -\frac{\partial H}{\partial q} \quad \text{a.e.} \\
H(q(t), p(t), u(t)) &= \max_{|v| \leq 1} H(q(t), p(t), v) \quad \text{a.e.}
\end{align}

Moreover, $M(q, p) = \max_{|v| \leq 1} H(q, p, v)$ is constant along $(q(\cdot), p(\cdot))$ and non-negative, with $H(q, p, u) = p \cdot (F + u G)$ the pseudo-Hamiltonian and $p$ is called the adjoint vector.

**Definition 2.2.** A triplet $(q(\cdot), p(\cdot), u(\cdot))$ solution of (2.3) and (2.4) is called an extremal. It is called regular if $u(t) = \text{sign}(p(t) \cdot G(q(t)))$ a.e. and bang-bang if it is regular and the number of switchings of $u(\cdot)$ is finite. An extremal is called singular if $p(\cdot) \cdot G(q(\cdot)) = 0$ everywhere. We denote by $\sigma_+$, $\sigma_-$ and $\sigma_s$ respectively bang with $u = +1$, $u = -1$ and singular extremals. Extremals satisfying the boundary conditions are called BC-extremals.

**Proposition 2.3.** If the control $u(\cdot)$ is singular on $[0, T]$ (for the extremity mapping), with $q(\cdot)$ the associated trajectory, then there exists $p(\cdot)$ such that $(q(\cdot), p(\cdot), u(\cdot))$ is a singular extremal.

2.2.2. Computation of singular trajectories. The Lie bracket of two $C^\infty$ vector fields $X, Y$ on $V$ is computed with the convention:

$$[X, Y](q) = \frac{\partial Y}{\partial q}(q) X(q) - \frac{\partial X}{\partial q}(q) Y(q),$$

and denoting $H_X, H_Y$ the Hamiltonian lifts: $H_X(z) = p \cdot X(q)$, $H_Y(z) = p \cdot Y(q)$, with $z = (q, p) \in V \times \mathbb{R}^n$, the Poisson bracket reads:

$$\{H_X, H_Y\} = \frac{\partial H_Y}{\partial p} \frac{\partial}{\partial q} - \frac{\partial H_X}{\partial q} \frac{\partial}{\partial p}.$$

Differentiating twice $p(\cdot) \cdot G(q(\cdot))$ with respect to the time $t$, one gets:

**Proposition 2.4.** Singular extremals $(z(\cdot), u(\cdot))$ are solutions of the following equations:

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

$$\{H_F, \{H_F, H_G\}\}(z(t)) + u(t) \{H_G, \{H_F, H_G\}\}(z(t)) = 0.$$

If $\{H_G, \{H_F, H_G\}\} \neq 0$ along the extremal, then the singular control is called of minimal order and it is given by the dynamic feedback:

$$u_s(z(t))) = -\frac{\{H_F, \{H_F, H_G\}\}(z(t))}{\{H_G, \{H_F, H_G\}\}(z(t))}.$$

From the above proposition one gets:

**Corollary 2.5.** If $u(\cdot) = 0$ is a singular control on $[0, T]$ then one has:

$$\text{ad}^k H_F \cdot H_G(z(t)) = p(\cdot) \cdot (\text{ad}^k F \cdot G(q(t))), \quad \forall k \geq 0,$$

with $\text{ad} F \cdot G = [F, G]$, $\text{ad} H_F \cdot H_G = \{H_F, H_G\}$.

The following lemma is useful [5, p. 213].

**Lemma 2.6.** For generic pair $(F, G)$ for the Whitney topology, $\dim \text{span}\{\text{ad}^k F \cdot G, k \geq 0\} = n.$
Corollary 2.7. Up to the time reparameterization
\[
\frac{ds}{dt} = \frac{1}{\{H_G, \{H_F, H_G\}\}(z(t))},
\]
singular extremals of minimal order are solutions of the analytic differential equation:
\[
\frac{dz}{ds} = X(z),
\]
with
\[
X = \left(\{H_G, \{H_F, H_G\}\} F - \{H_F, \{H_F, H_G\}\} G\right) \frac{\partial}{\partial q}
- \left(\{H_G, \{H_F, H_G\}\} \frac{\partial F}{\partial q} - \{H_F, \{H_F, H_G\}\} \frac{\partial G}{\partial q}\right) \frac{\partial}{\partial p},
\]
with two constraints \(H_F(z) = \{H_F, H_G\}(z) = 0\).

2.2.3. Classification of singular extremals and time optimality properties. In this section we recall results of [12] concerning singular extremals. We consider the \(C^\omega\)-single input control system relaxing the control bound \(|u| \leq 1\)
\[
\dot{q} = F(q) + uG(q), \quad u \in \mathbb{R}.
\]
Let \(\gamma(t) = (q(t), p(t)), \quad t \in [0, T]\) be a reference singular extremal of minimal order and we assume that \(t \mapsto q(t)\) is one-to-one. Assuming \(F, G\) not collinear along \(q(\cdot)\) one can assume that \(q(\cdot)\) is a singular extremal associated to \(u_\ast \equiv 0\). The first order Pontryagin cone \(K(t)\) is the subspace of codimension \(\geq 1\) generated by the vectors \(\text{ad}^k F \cdot G(q(t)), \quad k \geq 0\). We introduce the following generic assumptions
\[
\text{(H1) } \forall t \in [0, T], \quad \text{ad}^2 F \cdot G(q(t)) \notin K(t)
\]
\[
\text{(H2) } \forall t, K(t) \text{ is exactly of codimension one and generated by the vectors } \{\text{ad}^k F \cdot G(q(t)): \quad k = 0, \ldots, n - 2\}.
\]
\[
\text{(H3) If } n \geq 3, \quad \forall t \in [0, T], F(q(t)) \in \text{span}\{\text{ad}^k F \cdot G(q(t)): \quad k = 0, \ldots, n - 3\}.
\]
Under these assumptions, the problem is normal that is the adjoint vector \(p(\cdot)\) associated to \(q(\cdot)\) is unique up to a factor and \(\forall t \in [0, T], p(t)\) is orthogonal to \(K(t)\). Orienting \(p(\cdot)\) using the convention of the Maximum Principle: \(\langle p(t), F(q(t)) \rangle \geq 0\), the singular trajectory is called
- **Hyperbolic** if \(\langle p(t), \text{ad}^2 G \cdot F(q(t)) \rangle > 0\).
- **Elliptic** if \(\langle p(t), \text{ad}^2 G \cdot F(q(t)) \rangle < 0\).
- **Exceptional** if \(\langle p(t), F(q(t)) \rangle = 0\).

Note that the condition \(\langle p(t), \text{ad}^2 G \cdot F(q(t)) \rangle \geq 0\) amounts to the generalized Legendre-Clebsch condition
\[
\frac{\partial}{\partial u} \frac{d^2}{dt^2} \frac{\partial H}{\partial u}(\gamma(t)) \geq 0
\]
and according to the higher-order maximum principle [20], this condition is a necessary (small) time minimization condition. The key result is the following.

Theorem 2.8. Under assumptions \((H_1), (H_2)\) and \((H_3)\) an exceptional or hyperbolic (respectively elliptic) is time minimizing (respectively time maximizing) on \([0, T]\) with respect to all trajectories contained in a \(C^0\)-neighbourhood of \(q(\cdot)\) if \(T < t_{1c}\) where \(t_{1c} > 0\) is called the first conjugate time along \(q(\cdot)\).
Algorithms to compute the first conjugate times are described in [4] and are implemented in the HamPath code.

This gives a complete characterization of the time minimality status of the singular extremals under (generic) assumptions. The control bound $|u| \leq 1$ has been relaxed. In this classification, the control bound $|u_s| \leq 1$, that is the singular arc is admissible to complete the analysis. The case $|u_s| = 1$ corresponds to saturation. If $|u_s| > 1$, the singular control is not admissible and with some abuse, we shall use the terminology parabolic arc for the corresponding trajectory. See [21] for this terminology.

3. Time minimal saturation of a single spin

We refer to [34, 13] for the standard concepts of regular synthesis used in our analysis. We have two cases:

- case 1: fix the initial point to be the North Pole $N = (0,1)$;
- case 2: fix the final point to be the center $O = (0,0)$.

We shall only analyze the first case. We complete results from [11] and give new insights.

3.1. Lie brackets computations. The system (2.2) is written as:

$$\frac{dq}{dt} = F(q) + u G(q), \quad |u| \leq 1,$$

with

$$F(q) = -\Gamma y \frac{\partial}{\partial y} + \gamma (1 - z) \frac{\partial}{\partial z}, \quad G(q) = -z \frac{\partial}{\partial y} + y \frac{\partial}{\partial z},$$

and we can write

$$F(q) = Aq + a, \quad A = \begin{pmatrix} -\Gamma & 0 \\ 0 & -\gamma \end{pmatrix}, \quad a = \begin{pmatrix} 0 \\ \gamma \end{pmatrix}, \quad G(q) = Bq, \quad B = \begin{pmatrix} 0 & -1 \\ 1 & 0 \end{pmatrix}.$$

The system can be lifted on the semi-direct product $GL(2, \mathbb{R}) \ltimes \mathbb{R}^2$ acting on the $q$-space by the action $(A, a) \cdot q = Aq + a$, and where the Lie bracket rule is: $[(A, a), (B, b)] = ([A, B], Ab - Ba)$, with $[A, B] = AB - BA$ the commutator. One writes

$$A = \mu I_2 + \begin{pmatrix} \lambda & 0 \\ 0 & -\lambda \end{pmatrix},$$

with $\mu = -\frac{\Gamma + \gamma}{2}$, which is zero if and only if $\gamma = -\Gamma$ and $\lambda = \frac{\delta}{2}$, where $\delta = \gamma - \Gamma$.

The case $\delta = 0$ is the case of water species. Otherwise, we have:

**Lemma 3.1.** If $\delta \neq 0$, the Lie algebra generated by $(A, a), (B, 0)$ is $gl(2, \mathbb{R}) \oplus \mathbb{R}^2$.

This provides a rough classification of the problems between the simple water case and the general case.

Moreover, all the Lie brackets can be easily computed. They are listed next, up to length 4.

Length 1.

$$F(q) = -\Gamma y \frac{\partial}{\partial y} + \gamma (1 - z) \frac{\partial}{\partial z}, \quad G(q) = -z \frac{\partial}{\partial y} + y \frac{\partial}{\partial z},$$
The collinearity set \( \mathcal{C} \) is defined as the set where \( F \) and \( G \) are linearly dependent and the singular set \( \mathcal{S} \) is where \( F \) and \( [F,G] \) are collinear. Computing, one has:

**Lemma 3.2.** The collinearity set \( \mathcal{C} \) is given by \( y z (1 - z) - \Gamma y^2 = 0 \), thus \( O \) and \( N \) belong to \( \mathcal{C} \). Under the assumption \( 0 < y \leq 2 \Gamma \), \( \mathcal{C} \) is an ellipse contained in the Bloch ball. Besides, for each point \( q \) of \( \mathcal{C} \), except \( O \), there exists \( u \) such that \( q \) is an equilibrium point of the dynamics \( F + uG \).

The dynamics reads \( F(q) + u G(q) = Mq + a \), \( M = A + u B \), so for the maximal amplitude \( u = +1 \), the corresponding equilibrium point is

\[
O_1 = -M^{-1}a = \frac{\gamma}{u^2 + \gamma \Gamma} \begin{pmatrix} -u \\ -1 \end{pmatrix} = \frac{\gamma}{1 + \gamma \Gamma} \begin{pmatrix} -1 \\ 1 \end{pmatrix} \in \mathcal{C}
\]

and it is contained in the sector \( y < 0 \). We define in the same way \( O_{-1} \) for \( u = -1 \).

**Lemma 3.3.** The singular trajectories are located on \( \mathcal{S} \) which is given by \( y(\gamma - 2 \delta z) = 0 \). Hence, it is the union of the \( z \)-axis of revolution \( y = 0 \) and the horizontal line \( z = \gamma/(2 \delta) \), providing \( \delta \neq 0 \). Under the assumption \( 0 < \gamma \leq 2 \Gamma \), \( \mathcal{S} \) intersects the Bloch ball if and only if \( 3 \gamma \leq 2 \Gamma \). In this case, \( z_s = \gamma/(2 \delta) \in [-1,0) \).

The singular control is given by solving

\[
D'(q) + u D(q) = 0,
\]

with \( D = \det(G,[F,[F,G]]) \) and \( D' = \det(G,[F,[F,G]]) \). For \( y = 0 \), \( D(q) = -z(\gamma - 2 \delta z) \) and \( D' = 0 \). Hence, the singular control is zero and the singular trajectories are solution of \( \dot{y} = -\Gamma y = 0 \), \( \dot{z} = \gamma(1 - z) \). The North Pole is an equilibrium which is a stable node for \( 0 < \gamma \leq 2 \Gamma \). Along the horizontal singular line, i.e. for \( z = \gamma/(2 \delta) \), one has \( D(q) = -2 \delta y^2 \), \( D'(q) = \gamma y(2 \Gamma - \gamma) \) and the singular control denoted \( u_s \) is given by \( u_s(q) = \gamma(2 \Gamma - \gamma)/(2 \delta y) \). Hence, along the singular horizontal direction, the singular flow is:

\[
\dot{y} = -\Gamma y - \gamma^2(2 \Gamma - \gamma)/(4 \delta^2 y), \quad \dot{z} = 0,
\]

and one has \( u_s \to \pm \infty \) when \( y \to 0^\mp \).
Main assumptions. Under the physical assumptions
\[ 0 < \gamma \leq 2\Gamma, \]
the Bloch ball is invariant for the dynamics. Besides, under the physical assumption and
\[ |\delta| < 2, \]
any positive bang arc solution of \( \ddot{q} = F(q) + G(q) = Mq + a \) spirals around \( O_1 \) and converges to it (\( O_1 \) is a stable spiral), since in this case \( M \) has two complex conjugate eigenvalues with negative real part. Likewise, negative bang arcs spirals around \( O_{-1} \).

Proposition 3.4. If \( \gamma > 0 \) and \( |\delta| < 2 \), then the solution of \( \ddot{q} = F(q) + G(q) \), \( q(0) = q_0 \) is given by:
\[
q(t) = \exp((\alpha - \gamma)t) \left( \begin{array}{c} \frac{\alpha}{\beta} \sin(\beta t) + \cos(\beta t) \\ \frac{1}{\beta} \sin(\beta t) - \frac{\alpha}{\beta} \sin(\beta t) + \cos(\beta t) \end{array} \right) (q_0 - O_1) + O_1,
\]
where \( \alpha = \delta/2 \) and \( \beta = \sqrt{1 - \alpha^2} \). The solution is quasi-periodic of period \( T = 2\pi/\beta \).

Proposition 3.5. Let \((y(\cdot), z(\cdot))\), with associated control \( u(\cdot) \), be a trajectory solution of (2.2). Then, \((-y(\cdot), z(\cdot))\) with control \(-u(\cdot)\) is also solution of (2.2).

This discrete symmetry allows us to consider only trajectories in the domain \( y \leq 0 \) of the Bloch ball.

**Notation.** We denote by \( S_1 \) the intersection of the positive bang arc \( \sigma_+ \) issued from the North Pole with the horizontal singular line \( z = z_s = \gamma/2\delta \), and by \( S_1' \) the intersection with the vertical singular line \( y = 0 \). \( S_1 \) and \( S_1' \) may not exist for specific values of \((\gamma, \Gamma)\). We denote by \( S_3 \) the point on the horizontal singular line such that \( \sigma_+ \) is tangent to this line at this point, i.e. \( S_3 \) is a saturation point for the singular control \( u_s \) \((u_s(S_3) = 1)\), and by \( S_3' \) the intersection of the bang arc \( \sigma_+ \) issued from \( S_3 \) with the axis \( y = 0 \). If \( 0 < \gamma \leq 2\Gamma \) is satisfied, then \( S_3 \) is in the domain \( y < 0 \) if \( \delta < 0 \). We define in the same way \( S_{-3} \) for \( u = -1 \). See Fig. 1 to visualize on an example: the singular and collinearity sets, the points \( S_1, S_1', S_3, S_3' \)...

3.3. Switching function, the concept of bridge and the \( \theta \) function.

3.3.1. **Switching function.** Let \( z(t), t \in [0, T], \) be an extremal curve. The switching function is defined as \( \Phi(t) = p(t) \cdot G(q(t)) \). A time \( t \) is called an ordinary switching time if \( \Phi(t) = 0 \) and \( \dot{\Phi}(t) \neq 0 \), i.e. \( p(t) \cdot G(q(t)) = 0 \) and \( p(t) \cdot [F, G][q(t)] \neq 0 \). In the 2D-case, outside the collinearity set, one can write:
\[
[F, G](q) = \alpha(q) F(q) + \beta(q) G(q),
\]
with
\[
\alpha(q) = \frac{\det(G(q), [F, G](q))}{\det(G(q), F(q))}.
\]
At an ordinary switching time \( t \), one has:
\[
\mathrm{sign}(\dot{\Phi}(t)) = \mathrm{sign}(\alpha(q(t))),
\]
with the convention of the maximum principle, i.e. \( H_F \geq 0 \). If \( \alpha(q(t)) > 0 \), then we switch from an arc \( \sigma_- \) to an arc \( \sigma_+ \) and the converse if \( \alpha(q(t)) < 0 \). Fig. 2 gives the sign of \( \alpha \) inside the Bloch ball for fixed parameters \( \gamma \) and \( \Gamma \).
3.3.2. The concept of Bridge. An arc $\sigma_+$ or $\sigma_-$ corresponding to $u = +1$ or $u = -1$, is called a bridge on $[0, t]$ if the extremities correspond to non ordinary switching points, i.e. $\Phi(0) = \dot{\Phi}(0) = \Phi(t) = \dot{\Phi}(t) = 0$, see Fig. 3.

Remark 3.6. This concept is important and leads to a generalization in higher dimension, which plays an important role in the time minimal saturation of a pair of spins, but also in the contrast problem.
Figure 3. An extremal of the form $\sigma_+ \sigma_+ \sigma_+ \sigma_+$ is portrayed in the cotangent space, where $\sigma_+$ and $\sigma_-$ represent respectively bang and singular extremals. The second bang arc is a bridge.

Figure 4. The points $S_1$, $S_2$ and $S_3$ with different orders. On the left sub-graph: $(\gamma, \Gamma) = (0.1, 0.5)$ and on the right: $(\gamma, \Gamma) = (0.163, 0.5)$.

If a bridge $\sigma_+$, in the domain $y \leq 0$, connecting the horizontal and vertical singular lines exists, then we denote by $S_2$ the extremity on $z = \gamma/2\delta$, and $S'_2$ the other extremity on $y = 0$. We observe two cases when $S_2$ exists which are crucial for the analysis of the time minimal saturation problem of a single spin, see Fig. 4.

3.3.3. Analysis of two consecutive switching times. In the 2D-case, in order to analyze switchings, one proceeds as follows. Assume $0$ and $t$ be two consecutive switching times on an arc $\sigma_+$ or $\sigma_-$. Let $z(\cdot) = (q(\cdot), p(\cdot))$ denote the associated extremal. We have:

$$p(0) \cdot G(q(0)) = p(t) \cdot G(q(t)) = 0.$$  

We denote by $v(\cdot)$ the solution on $[0, t]$ of the variational equation

$$\dot{v}(t) = \left( \frac{\partial F}{\partial q}(q(t)) + u \frac{\partial G}{\partial q}(q(t)) \right) v(t), \quad v(t) = G(q(t)), \quad u = \pm 1.$$
This equation is integrated backwards from time \( t \) to 0. By construction, \( p(\cdot) \cdot v(\cdot) \) is constant and equal to zero. At time 0, one has \( p(0) \cdot v(0) = p(0) \cdot G(q(0)) = 0 \). Hence, \( p(0) \) is orthogonal to \( v(0) \) and to \( G(q(0)) \). Therefore, \( v(0) \) and \( G(q(0)) \) are collinear (\( p(\cdot) \) does not vanish, according to proposition 2.1). We introduce naturally the \( \theta(t) \) function which gives the angle between \( G(q(0)) \) and \( v(0) \) measured counterclockwise. One deduces that switchings occur at times 0 and \( t \) if

\[
\theta(t) = 0 \mod \pi
\]

and it can be tested using \( \det(G(q(0)),v(0)) = 0 \). This test may be used by contraposition in order to eliminate the possibility that bang-bang trajectories with at least two switching times (or more) are optimal. We have by definition

\[
v(0) = e^{-t \operatorname{ad}(F+uG)}(G(q(t))), \quad u = \pm 1,
\]

and in the analytic case, the \( \operatorname{ad} \)-formula gives:

\[
v(0) = \sum_{n \geq 0} \frac{(t)^n}{n!} \operatorname{ad}^n(F + uG) \cdot G(q(t)).
\]

The computation can be made explicit on a Lie group since determining \( \exp(t \operatorname{ad}(F+uG)) \) amounts to compute a Jordan form of the linear operator \( \operatorname{ad}(F+uG) \) defined by the Lie brackets.

### 3.4. Optimality status

A first step in the optimality analysis is to discriminate between small time minimizing or maximizing singular trajectories using the high-order maximum principle and Theorem 2.8. In dimension 2, we introduce \( D'' = \det(G,F) \) and \( D'' = 0 \) is the collinearity set \( \mathcal{C} \). Singular lines are fast if \( DD'' > 0 \) and slow if \( DD'' < 0 \). Using these conditions, if \( 0 < \gamma \) and \( \delta < 0 \), then the horizontal singular line \( z = z_s \) is fast if \( y \neq 0 \). For the interesting case when the horizontal singular line cuts the Bloch ball, \( i.e. \) when \( 0 < 3\gamma \leq 2\Gamma \), the singular horizontal line is fast. Moreover, it is parabolic on \( (S_3, S_{-3}) = \{ (1 - \lambda)S_3 + \lambda S_{-3} \mid \lambda \in (0,1) \} \) when \( y \neq 0 \) and hyperbolic otherwise (except of course at \( S_3 \) and \( S_{-3} \)). On the other hand, assuming \( 0 < \gamma \leq 2\Gamma \) and considering only the interesting part inside the Bloch ball, then we have the following: if \( \delta < 0 \), then the vertical line is hyperbolic for \( z_s < z < 1 \) and elliptic for \( -1 \leq z < z_s \). If \( \delta > 0 \), then it is hyperbolic for \( -1 \leq z < 1 \).

Theorem 2.8 combined with the generalized Legendre-Clebsch condition gives information about the local optimality of the singular extremals. In the 2D-case, global optimality can be analyzed using the clock form \( \omega = p \cdot dq \) with \( p \cdot G = 0 \) and \( p \cdot F = 1 \). The clock form may be used to compare two trajectories with the same extremities whenever they do not cross the collinearity set. In the time minimal saturation problem of a single spin, the collinearity set plays a crucial role and we must use the \( \theta \) function defined in section 3.3 to eliminate the possibility to have two consecutive ordinary switching times.

To complete the study, we have to analyze the behavior of optimal trajectories in the neighborhood of some particular points, named Frame Points corresponding to isolated singularities: intersection of the collinearity locus with singular locus or singular loci [13]. In the time minimal saturation problem, there exists two phenomena to analyze. The horizontal singular line being admissible up to a saturation
point $S_3$, there is a birth of a switching locus $\Sigma_3$ connecting the horizontal and vertical singular lines. This is related to the concept of bridge and this phenomenon is referred as the SiSi singularity [11]. The time minimal synthesis, with initial point $N$, is represented on Figs. 9, 10 and 11. The interaction between the collinearity and singular sets near the North Pole is the second phenomenon to analyze and is referred as the SiCo singularity [11]. Near the North Pole, only bang-bang trajectories with at most one switching are optimal (see the top part of Fig. 9).

According to the sections 3.5, 3.6, 3.7 and 3.8 and according to the Figs. 9, 10 and 11, we have the following results.

**Theorem 3.7.** Let us denote by $\sigma_N^+$ the positive bang arc starting from the North Pole with $S_1, S'_1$ respectively the intersection points (they may not exist) with the horizontal, vertical singular line. Let us denote by $\sigma_b^+$ the bridge with $S_2$ and $S'_2$ as extremities, by $\sigma_h^+, \sigma_v^+$ respectively a horizontal, vertical singular arc. Let $S_3$ denote the saturation point on the horizontal singular line and $S'_3$ the intersection of the bang arc starting from $S_3$ with the axis $y = 0$.

For parameters $(\Gamma, \gamma)$ satisfying the physical constraints $0 < \gamma \leq 2\Gamma$, and such that the points $S'_1$ and $S'_3$ are below the origin $O$, then the minimal time trajectory to steer the spin from $N$ to $O$ is $NS_1S_2S'_2O$, i.e. it is of the form $\sigma_N^+\sigma_h^+\sigma_b^+\sigma_v^+$ (no empty arcs), if $\sigma_N^+$ intersects the horizontal singular line strictly before $S_2$, that is $S_1 < S_2$. Otherwise, the optimal trajectory is $NS'_1O$, i.e. of the form $\sigma_N^+\sigma_v^+$.

**3.5. Parameters and practical cases.** The optimal syntheses depend on the parameters $(\gamma, \Gamma)$. We define in this section the domain of interest of the parameters and partition it in four main sub-domains which are denoted $A_1, A_2, B$ and $C$. This partitioning may be visualized on Fig. 5.

We have already seen that we assume $0 < \gamma \leq 2\Gamma$ and $|\delta| < 2$, with $\delta = \gamma - \Gamma$. Additionally, we require that the origin $O$ is accessible from the North Pole $N$, that is we impose that $S'_1$ is below $O$. The parameters such that $S'_1 = O$ are given by the following proposition:

**Proposition 3.8.** For $\gamma > 0$, we have:

$$S'_1 = O \iff \exp \left( (t_0 \beta + \pi) \frac{\alpha - \gamma}{\beta} \right) - \gamma = 0,$$

with

$$t_0 = \begin{cases} \frac{1}{\beta} \arctan \left( -\frac{\delta}{\alpha} \right) & \text{if } \delta < 0, \\ \frac{\pi}{2} & \text{if } \delta = 0, \\ \frac{1}{\beta} \left( \arctan \left( -\frac{\delta}{\alpha} \right) + \pi \right) & \text{if } \delta > 0. \end{cases}$$

Besides, to simplify the presentation, we restrict the analysis to the case where $S'_3$ is below $O$. This implies that the origin may be reached from the horizontal singular locus going through the bridge. We have the following relation on the parameters to impose $S'_3 = O$:

**Proposition 3.9.**

$$S'_3 = O \iff (2\Gamma^2 - \gamma \Gamma + 1) \exp \left( (\alpha - \gamma)t_0 \right) - 2|\delta| = 0.$$

---

1. We denote by $\Sigma$ the switching locus with strata $\Sigma_1, \Sigma_2, \ldots$
One can notice according to Fig. 5 that if \(0 < \gamma \leq 2\Gamma\) is satisfied and if the points \(S_1'\) and \(S_3'\) are below the origin \(O\), then \(|\delta| < 2\) holds. We restrict the analysis of the optimal syntheses to these values of parameters. Let us explain now how we partition this domain. We denote by \(A_1\) the sub-domain such that \(2\Gamma/3 \leq \gamma \leq 2\Gamma\) and \(S_1'\) is below \(O\). In this case, \(S_3'\) is necessarily below \(O\), the horizontal singular line (at \(z = z_s = \gamma/2\delta\)) does not cut the interior of the Bloch ball and the optimal synthesis is simple, see section 3.8. Now, when the horizontal singular line cuts the interior of the Bloch ball, that is for \(0 < \gamma < 2\Gamma/3\), this part of the singular locus does not play any role if the bang arc starting from the North Pole does not intersect it, that is if \(S_1\) does not exist. The limit case is when \(S_1 = S_3\), where \(S_3\) is the saturation point on the horizontal singular line. We denote by \(A_2\) this sub-domain of parameters and we have the same optimal syntheses in \(A_1\) and \(A_2\), see again section 3.8.

**Proposition 3.10.**

\[
S_1 = S_3 \iff (2\Gamma^2 - \gamma \Gamma + 1) \frac{\gamma}{2\delta} + \exp\left((\alpha - \gamma) \frac{\pi}{\beta}\right) = 0.
\]

The remaining part, when \(S_1\) exists, may be split in two. Either, \(S_1\) is before \(S_2\) (\(S_2\) is one of the extremities of the bridge) or after. We denote by \(B\) the sub-domain where \(S_2 < S_1\) (i.e. \(S_2\) strictly before \(S_1\)) and \(C\) where \(S_1 \leq S_2\). One can visualize the sub-domains \(A_1, A_2, B\) and \(C\) on the Fig. 5.

**Remark 3.11.** Each specie to study is characterized by its relaxation times \(T_1\) and \(T_2\), see Table 1. However, the main parameter is the ratio \(T_2/T_1\). According to Fig. 6, one can see that for any couple \((T_1, T_2)\) such that \(0 < T_2 < 2T_1/3\), then there exists \(\omega_{\text{max}} > 0\) such that the associated parameters \((\gamma, \Gamma)\), see eq. (2.1), belongs to the sub-domain \(C\). Hence, for any couple \((T_1, T_2)\) satisfying the physical constraint \(0 < T_2 < 2T_1\), there exists \(\omega_{\text{max}} > 0\) such that the associated parameters \((\gamma, \Gamma)\) belongs either to \(A_1\) or \(C\).
Figure 6. The sub-domains in $(\Gamma,\gamma/\Gamma)$ coordinates. One can notice that when $\Gamma$ tends to 0, then the slopes of the curves $S_1 = S_2$ and $S_1 = S_3$, in $(\Gamma,\gamma)$ coordinates, are equal to $2/3$.

Figure 7. The associated $(\gamma, \Gamma)$ parameters of cases from Table 1 with $\omega_{\text{max}} = 2\pi \times 32.3$ Hz. Except the water case which is in the sub-domain $A_1$, all the others cases belong to the sub-domain $C$.

In Fig. 7, we show for the practical cases from Table 1, the associated $(\gamma, \Gamma)$ parameters with $\omega_{\text{max}} = 2\pi \times 32.3$ Hz. Note that in the experiments, $\omega_{\text{max}}$ may be chosen up to 15 000 Hz but we consider here the same value as in [11]. The water case belongs to the domain $A_1$ while all the others cases are contained in $C$. In Fig. 8 is represented the slope $T_2/T_1 = \gamma/\Gamma$ with the particular case when $\omega_{\text{max}} = 2\pi \times 32.3$ Hz. One can notice that the fat case is the only one which crosses the sub-domains $A_2$, $B$ and $C$ but for the value $\omega_{\text{max}} = 2\pi \times 32.3$ Hz, it belongs to $C$.

3.6. The optimal synthesis in the sub-domain $C$. For the optimal synthesis in the sub-domain $C$, we have the following: the parameters satisfy $0 < 3\gamma < 2\Gamma$, the point $S'_3$ is below the origin $O$ and there exist $S_1$ and $S_2$ such that $S_1 \leq S_2 < S_3$, i.e. we are in the situation of the left sub-graph of Fig. 4. In this case, the horizontal singular line cuts the Bloch ball in the domain $-1 < z < 0$ and the global synthesis
is similar as the one presented in [11]. It is obtained gluing together the SiSi and SiCo cases and it is represented on the Fig. 9.

**Remark 3.12.** In [11], to obtain this optimal synthesis, it is assumed that $\omega_{\text{max}}$ is large enough. Remark 3.11 explains why this assumption is correct. However, this assumption is replaced here by geometric relations on the points $S_1$, $S_2$ and $S_3$.

The switching locus is formed by the positive bang arc starting from the North Pole (denoted $\sigma_N$) and reaching the horizontal singular arc at $S_1$ (it is denoted $\Sigma_1$ in the figure), by the horizontal singular segment $\Sigma_2$ between the points $S_1$ and $S_3$, the switching locus $\Sigma_3$ due to the saturation phenomenon and by the part of the vertical singular direction between $S'_2$ and $O$ (the $\Sigma_4$ segment), $S'_2$ being the extremity of the bridge on $y = 0$. The bang arc with $u = -1$ starting from $S_1$ splits the domain in two sub-domains, one with a bang-bang policy and the other containing a non trivial singular arc.

We have, as a corollary, that in this case, the optimal strategy to steer the system from the North Pole to the origin in minimum time is of the form $\sigma_N^r \sigma_s^h \sigma_s^v \sigma_s^v$, where $\sigma_s^h$ and $\sigma_s^v$ denote respectively horizontal and vertical singular arcs, and where $\sigma_s^v$ is the bridge.

**Remark 3.13.** Note that the switching locus has a complex structure, but due to the symmetry, all the cut points, i.e. the first points where the extremal trajectories cease to be optimal, are on the vertical $z$-axis where two symmetric solutions starting respectively on the left and right part of the Bloch ball intersect at the same time.

### 3.7. The optimal synthesis in the sub-domains $B$

For the optimal synthesis in the sub-domain $B$, we have the following: the parameters satisfy $0 < 3\gamma < 2\Gamma$, the point $S'_3$ is below the origin $O$ and there exist $S_1$ and $S_2$ such that $S_2 < S_1 \leq S_3$, i.e. we are in the situation of the right sub-graph of Fig. 4. In this case, the horizontal singular line cuts the Bloch ball in the domain $-1 < z < 0$ and still plays a role.
Figure 9. Schematic time minimal synthesis to steer a single spin system from the North Pole N to any point of the Bloch ball in the reachable set, for parameters \((\Gamma, \gamma) \in C\). An arbitrary zoom has been used to construct the figure. The set of \(\Sigma_i\) forms the switching surface \(\Sigma\) dividing the \(+1\) and \(-1\) areas respectively in red and blue. The minimal time trajectory to steer the spin from N to O is \(NS_1S_2S_3O\), i.e. it is of the form \(\sigma^N_+\sigma^h_+\sigma^v_+\sigma^-\) with horizontal \(\sigma^h\) and vertical \(\sigma^v\) singular arcs. The spin leaves the horizontal singular arc before the point \(S_3\) (where the control saturates the constraint) producing a bridge \(\sigma^b_+\) to reach the vertical singular line.

The switching locus is formed by the positive bang arc starting from the North Pole (denoted \(\sigma^N_+\)) and reaching the horizontal singular arc at \(S_1\) (denoted \(\Sigma_1\)), by the horizontal singular segment \(\Sigma_2\) between the points \(S_1\) and \(S_3\), by the switching locus \(\Sigma_3\) due to the saturation phenomenon from \(S_3\) to the intersection with \(\sigma^N_+\) (denoted \(S_1''\)), by the part of \(\sigma^N_+\) between \(S_1''\) and \(S_1\) (denoted \(\Sigma_3\)) and by the part of the vertical singular direction between \(S_1\) and O (the \(\Sigma_4\) segment), \(S_1''\) being the extremity of \(\sigma^N_+\) on \(y = 0\).

We have, as a corollary, that in this case, the optimal strategy to steer the system from the North Pole to the origin in minimum time is of the form \(\sigma^N_+\sigma^v_+\).

Remark 3.14. Note that the switching locus is located on the vertical \(z\)-axis due to the symmetry.

3.8. The optimal synthesis in the sub-domains \(A_1\) and \(A_2\). For the optimal synthesis in the sub-domains \(A_1\) and \(A_2\), we have the following: the parameters satisfy \(0 < \gamma \leq 2\Gamma\), the point \(S_1\) is below the origin O and \(\sigma^N_+\) does not intersect the horizontal singular line \(z = z_s = \gamma/2\delta\), that is \(S_1\) does not exist. In this case, the horizontal singular line does not play any role.
The switching locus is formed by the positive bang arc starting from the North Pole (denoted $\sigma_N^+$) and reaching the vertical singular arc at $S_1^+$ (denoted $\Sigma_1$) and by the part of the vertical singular direction between $S_1^+$ and $O$ (the $\Sigma_4$ segment).

We have, as a corollary, that in this case, the optimal strategy to steer the system from the North Pole to the origin in minimum time is of the form $\sigma_N^+ \sigma^v_s$ as in the sub-domain $B$.

**Remark 3.15.** Note that the switching locus is located on the vertical $z$-axis due to the symmetry.

### 3.9. Geometric comments and construction of a normal form.

**Lie brackets computations.** At the intersection $q_0$, of the horizontal and vertical lines the Lie algebraic structure up to order four is described by:
Figure 11. Schematic time minimal synthesis to steer a single spin system from the North Pole $N$ to any point of the Bloch ball in the reachable set, for parameters $(\Gamma, \gamma) \in A_2$. For $(\Gamma, \gamma) \in A_1$, the synthesis is the same but there is no horizontal singular line inside the Bloch ball. An arbitrary zoom has been used to construct the figure. The minimal time trajectory to steer the spin from $N$ to $O$ is $NS'_1O$, i.e. it is of the form $\sigma_N^\Sigma \sigma_s^\Sigma$.

This led to the algebraic characterization of some specific bridges, related to this case.

**Normal form.** A normal form to make an explicit evaluation of the synthesis in the neighbourhood of zero of size $\varepsilon$ is obtained by choosing parameters in such a way that the points $S_2$ and $S_3$ are contained in this neighbourhood.
Limit syntheses. Since the singular control \( u_s \) is \( L^1 \) when \( q \to q_0 \), stay on the horizontal singular line, one has:

**Proposition 3.16.** Relaxing \(|u| \leq 1\), we have two limit time minimal solutions to steer \( N \) to \( O \)

- \( 2\Gamma > 3\gamma > 0 : R\sigma_s^h \sigma_s^v \)
- \( 2\Gamma \leq 3\gamma : R\sigma_s^v \)

where \( R \) is the rotation with respect to the center \( O \) to reach either the horizontal singular line or the vertical singular line.

### 4. Time minimal saturation of a pair of spin-1/2 particles

#### 4.1. The model.

Let us consider a couple of spins with the same characteristics, i.e. the same relaxation times \( T_1 \) and \( T_2 \), but for which for each, the control field has different intensities, because of inhomogeneities. The system we consider is the following:

\[
\begin{align*}
\dot{q}_1(t) &= F(q_1(t)) + u(t) G(q_1(t)), \\
\dot{q}_2(t) &= F(q_2(t)) + u(t) (1 - \varepsilon) G(q_2(t)),
\end{align*}
\]

where \( q_i = (y_i, z_i), \ i = 1, 2, \) denote the coordinates of each system and where the vector fields \( F \) and \( G \) are given by eq. (2.2), see sections 2.1 and 3.1. The term \((1 - \varepsilon), \varepsilon > 0 \) small, is the rescaling factor of the control maximal amplitude. We define the time minimal saturation problem of a pair of spin-1/2 particles as the following affine control problem with Mayer cost:

\[
(P_{BS}) \begin{cases}
 J(u(\cdot), t_f) = t_f \to \min \\
 \dot{q}(t) = F(q(t)) + u(t) G(q(t)), \ & |u(t)| \leq 1, \ t \in [0, t_f], \ q(0) = q_0, \\
 q(t_f) = q_f,
\end{cases}
\]

where \( q = (q_1, q_2) = (y_1, z_1, y_2, z_2), q_0 = (0, 1, 0, 1), q_f = (0, 0, 0, 0), \) and where we use the notation

\[
\begin{align*}
F(q) &= F(q_1) \frac{\partial}{\partial q_1} + F(q_2) \frac{\partial}{\partial q_2}, \\
G(q) &= G(q_1) \frac{\partial}{\partial q_1} + (1 - \varepsilon) G(q_2) \frac{\partial}{\partial q_2}.
\end{align*}
\]

We have the following Lie brackets up to order 3:

\[
\begin{align*}
[F, G](q) &= [F, G](q_1) \frac{\partial}{\partial q_1} + (1 - \varepsilon) [F, G](q_2) \frac{\partial}{\partial q_2}, \\
[F, [F, G]](q) &= [F, [F, G]](q_1) \frac{\partial}{\partial q_1} + (1 - \varepsilon) [F, [F, G]](q_2) \frac{\partial}{\partial q_2}, \\
[G, [F, G]](q) &= [G, [F, G]](q_1) \frac{\partial}{\partial q_1} + (1 - \varepsilon)^2 [G, [F, G]](q_2) \frac{\partial}{\partial q_2}.
\end{align*}
\]
4.2. Some prior theoretical results. We present some theoretical results in relation with numerical homotopies. They are based on Lie brackets computations.

**Lemma 4.1.** \( \text{span}\{\text{ad}^k F \cdot G; \ k \geq 0\} = \text{span}\{\text{ad}^k F \cdot G; \ k = 0, 1, 2, 3\}. \)

**Proposition 4.2.** For all \( q \) in the Bloch ball, for all pair of parameters \((\Gamma, \gamma)\), we have \( \det(G(q), [F,G](q), [F,[F,G]](q), [F,[F,[F,G]]](q)) = 0 \) and every solution of \( F(q) \) is a (smooth) singular trajectory.

**Lemma 4.3.** In the water case \( \gamma = \Gamma \), \( \text{ad}^2 F \cdot G \) is constant and collinear to \([F,G]\).

**Proposition 4.4.** In the water case \( \gamma = \Gamma \), the only singular trajectories are the (smooth) solutions of \( F(q) \).

Next, we present some accessibility criteria.

**Lemma 4.5.** A necessary condition to steer the North pole \((0,1,0,1)\) to \((0,0,0,0)\) is that for the second spin the arc \( \sigma_+ \) starting from \((0,1)\) reaches the vertical axis \( y_2 = 0 \) at a point \((0,z), \ z \leq 0\).

**Definition 4.6.** The center \( O \) is called \( \eta \)-reachable if there exists \( O' \) of the collinearity locus such that \( |O - O'| \leq \eta \) and \( O' \) is a point of the collinear locus such \( (F + u_0 G)(O') = O \) and \( O' \) is globally stable.

4.3. First numerical results and validation with LMI methods.

4.3.1. Direct approach (Bocop). The Bocop software [3] implements a so-called direct transcription approach, where the continuous optimal control problem (OCP) is transformed into a nonlinear programming (NLP). The reformulation is done by a discretization of the time interval, with an approximation of the dynamics of the system by a generalized Runge-Kutta scheme. This is part of direct local collocation methods. We refer the reader to for instance [2], [18] and [30] for more details on direct transcription methods and NLP algorithms.

We present local solutions obtained by the Bocop software for the bi-saturation problem \( (P_{BS}) \) in the Deoxygenated blood case (denoted \( C_1 \)), the Oxygenated blood case \( (C_2) \) and the Cerebrospinal fluid case \( (C_3) \) with \( \omega_{\text{max}} = 2\pi \times 32.3 \) and the Water case \( (C_4) \) with a larger value of \( \omega_{\text{max}} \). Let us mention that for \( C_4 \), the value of \( \omega_{\text{max}} \) is larger than \( 2\pi \times 32.3 \) just to obtain a control law with longer bang arcs, to visualize better the solution. The parameter \( \varepsilon \) is fixed to 0.1 and let us recall that the associated relaxation times are given in Table 1. The time evolution of the state variables \( q = (q_1, q_2) \) and of the control variable \( u \) are represented in Figs. 12–15, the optimal time is given in Table 2 and is compared with the optimal time for the saturation of the single spin. Note that for each case, there is one more Bang-Singular sequence (the first one) in the bi-saturation problem than in the mono-saturation problem. Besides, the remaining part looks like the strategy to steer one single spin to the center of the Bloch ball except that for \( C_1, C_2 \) and \( C_3 \), the penultimate singular part of the trajectory do not follow exactly the horizontal lines \( z_1 = z_2 = z_3 = \gamma/2\delta, \ \delta = \gamma - \Gamma \). Finally, note that both spins reach the vertical line \( y_1 = y_2 = 0 \) with \( z_1 = z_2 \) before the final singular arc with \( u = 0 \) to reach the center of the Bloch ball at the same time. In other words, the spins are synchronised at this time.
techniques provide for the saturation problem, lower bounds on the global optimal problem are globally optimal using moment/LMI techniques. More precisely, these techniques provide for the saturation problem, lower bounds on the global optimal

<table>
<thead>
<tr>
<th>Case</th>
<th>$\Gamma$</th>
<th>$\gamma$</th>
<th>$t_f$ (2 spins)</th>
<th>$t_f$ (1 spin)</th>
</tr>
</thead>
<tbody>
<tr>
<td>$C_1$</td>
<td>$9.855 \times 10^{-2}$</td>
<td>$3.65 \times 10^{-3}$</td>
<td>44.769</td>
<td>42.685</td>
</tr>
<tr>
<td>$C_2$</td>
<td>$2.464 \times 10^{-2}$</td>
<td>$3.65 \times 10^{-3}$</td>
<td>113.86</td>
<td>110.44</td>
</tr>
<tr>
<td>$C_3$</td>
<td>$1.642 \times 10^{-2}$</td>
<td>$2.464 \times 10^{-3}$</td>
<td>168.32</td>
<td>164.46</td>
</tr>
<tr>
<td>$C_4$</td>
<td>$9.855 \times 10^{-2}$</td>
<td>$9.855 \times 10^{-2}$</td>
<td>15.0237</td>
<td>8.7445</td>
</tr>
</tbody>
</table>

Table 2. Cases treated numerically corresponding respectively to the Desoxygenated case ($C_1$), the Oxygenated case ($C_2$), the Cerebrospinal fluid case ($C_3$) and the Water case ($C_4$). The 5th (resp. 4th) column gives the final time found by Bocop for the saturation of one spin (resp. two spins with $B_1$-inhomogeneity). The parameter $\varepsilon$ is fixed to 0.1.

**Figure 12.** Desoxygenated blood case ($C_1$) with RF-inhomogeneity ($\varepsilon = 0.1$). Trajectories for spin 1 and 2 in the ($y,z$)-plane are portrayed in the first two subgraphs. The corresponding control is drawn in the right subgraph. The horizontal lines $z_1 = z_2 = z_s = \gamma/2\delta$, $\delta = \gamma - \Gamma$, is represented by dashed lines. Note that the last bang arc is not well captured by the direct solver because it is too short.

**Figure 13.** Oxygenated blood case ($C_2$) with RF-inhomogeneity ($\varepsilon = 0.1$). Trajectories for spin 1 and 2 in the ($y,z$)-plane are portrayed in the first two subgraphs. The corresponding control is drawn in the right subgraph. The horizontal lines $z_1 = z_2 = z_s = \gamma/2\delta$, $\delta = \gamma - \Gamma$, is represented by dashed lines. Note that the last bang arc is not well captured by the direct solver because it is too short.

4.3.2. LMI method (*GloptiPoly*). A crucial step is to check whether the local optimal times presented in Table 2 and obtained with Bocop for the saturation problem are globally optimal using moment/LMI techniques. More precisely, these techniques provide for the saturation problem, lower bounds on the global optimal
time which can be used as a validation of the global optimality if the gap between
the lower bound obtained from moment/LMI techniques and the time obtained
by the direct method is small. This combination of techniques has already been
successful in the contrast problem [8] by nuclear magnetic resonance in medical
imaging.

The moment approach is a global optimization method which relaxes a non linear
optimal control problem using measures as a linear programming (LP) problem. In
the case where the data are polynomials, we can handle these measures by their
moment sequences. Using powerful certificate coming from algebraic geometry, e.g.
Putinar’s Positivstellensatz [32], this leads to an infinite dimensional LMI problem
which can be truncated to a finite set of moments. The sequence of optimal values
associated to these truncated problems converges to the optimal value of problem
(\(P_{BS}\)), that we denote by \(T_{\text{min}}^*\). We present a classical formulation and an alternative
one which exploits the structure of the problem, based on [16]. Note that we present
the method on the bi-saturation problem but it is straightforward to adapt the
explanations for the mono-saturation problem.

Notations. \(B_n\) is the unit ball of dimension \(n\), \(\mathcal{M}^+(Z)\) is the set of finite, positive
Borel measures supported on compact set \(Z\) and \(\int f(z) \, d\mu\) denotes the integration
of a continuous function \(f \in C(Z)\) with respect to \(\mu \in \mathcal{M}^+(Z)\).

Step 1: linear program on measures.
• 1st formulation: Moment approach using occupation measures.

Following [25], the problem \((P_{BS})\) can be embed into the linear program on measures:

\[
T_{LP} = \inf_{\mu, \mu_f} \int \! d\mu_f \quad \text{s.t.} \quad \int \left( \frac{\partial v}{\partial t} + \frac{\partial v}{\partial q} (F + uG) \right) \, d\mu = \int v(\cdot, q_f) \, d\mu_f - v(q_0), \quad \forall v \in \mathbb{R}[t, q],
\]

\[
\mu \in \mathcal{M}^+([0, T] \times Q \times U), \quad \mu_f \in \mathcal{M}^+(Q_f),
\]

and where \(T\) is fixed, \(Q_f = [0, T] \times \{(0,0,0,0)\}\), \(Q = B_2 \times B_2\) are admissable state sets and \(U = B_1\) is the admissible control set. Given any admissible pair \((q(\cdot), u(\cdot))\) for \((P_{BS})\), it corresponds a measure \(\mu\) admissible for (4.1) achieving the same cost, hence \(T_{\text{min}}^* \geq T_{LP}\). Moreover, according to Theorem 3.6 (ii) of [25], there is no optimality gap and

\[
T_{\text{min}}^* = T_{LP}.
\]

**Remark 4.7.** Since the dynamic is autonomous, the time variable can be removed and the LP problem becomes

\[
T_{LP} = \inf_{\mu} \int \! d\mu \quad \text{s.t.} \quad \int \frac{\partial v}{\partial q} (F + uG) \, d\mu = v(q_f) - v(q_0), \quad \forall v \in \mathbb{R}[q],
\]

\[
\mu \in \mathcal{M}^+(Q \times U).
\]

• 2nd formulation: Moment approach using modal occupation measures.

In the first formulation (4.1) (respectively (4.2)), measures are supported on the set \([0, T] \times Q \times U\) (respectively \(Q \times U\)) of dimension \(1 + 4 + 1 = 6\) (respectively \(5\)) and we expect them to be located on the optimal trajectory \((q^*(\cdot), u^*(\cdot))\). An alternative formulation [16] is to model controls by measures such that the measures are supported on \(Q\) only. Indeed, note that the dynamic in \((P_{BS})\) is affine in the control \(u\) which takes its values inside the polytope \(\text{conv}\{-1, +1\}\). This optimal control problem can be written as a switching system with two modes, the first mode corresponding to \(u = +1\) and the second mode corresponding to \(u = -1\). This leads to consider

\[
T_{LP'} = \inf_{\mu_1, \mu_2, \mu_f} \int \! d\mu_f \quad \text{s.t.} \quad \forall v \in \mathbb{R}[t, q], \quad \int \left( \frac{\partial v}{\partial t} + \frac{\partial v}{\partial q} (F + G) \right) \, d\mu_1 + \int \left( \frac{\partial v}{\partial t} + \frac{\partial v}{\partial q} (F - G) \right) \, d\mu_2 = \int v(\cdot, q_f) \, d\mu_f - v(0, q_0),
\]

\[
\mu_1, \mu_2 \in \mathcal{M}^+([0, T] \times Q), \quad \mu_f \in \mathcal{M}^+(Q_f)
\]
where $T$ is fixed, $Q_f = [0, T] \times \{(0, 0, 0, 0)\}$ and $Q = B_2 \times B_2$ are admissible state sets.

**Step 2: Moment SDP.** An important feature of the problems (4.1)-(4.3) is their algebraic structure: the dynamic is polynomial and the sets $Q$ and $U$ are compact basic semi-algebraic sets. In these settings, it is possible to handle the measures by their moments which leads to a semi-definite program on countably many moments. Let us introduce for a multi-index $\alpha = (\alpha_1, \ldots, \alpha_p) \in \mathbb{N}^p$ and $y = (y_1, \ldots, y_p) \in \mathbb{R}^p$, the notation $|\alpha|_1 = \sum_{i=1}^{p} \alpha_i$ and $y^\alpha$ which stands for the monomial $y_1^{\alpha_1} \cdots y_p^{\alpha_p}$. Then, we denote by $\mathbb{N}^p_d$ the set $\{\alpha \in \mathbb{N}^p \mid |\alpha|_1 \leq d\}$.

**Definition 4.8.** The moment of order $\alpha \in \mathbb{N}^p$ of a measure $\mu$ supported on $Z \subset \mathbb{R}^p$ is the real $y_\alpha = \int z^\alpha \, d\mu$. Besides, $\mu \in \mathcal{M}(z)$ is said to be a representing measure for a sequence $(y_\alpha)_\alpha$ if $y_\alpha = \int z^\alpha \, d\mu$ for all $\alpha \in \mathbb{N}^p$.

**Definition 4.9.** Given an arbitrary sequence of reals $(y_\alpha)_\alpha$, we define the Riesz linear functional $l_y : \mathbb{R}[z] \to \mathbb{R}$ by $l_y(z^\alpha) = y_\alpha$ for all $\alpha \in \mathbb{N}^p$.

**Definition 4.10.** The moment matrix $M_d(y)$ of order $d$ is such that $l_y(p(z)^2) = p'M_d(y)p$ for all polynomials $p(z)$ of degree $d$ whose coefficients are denoted by the vector $p$. In particular, the $(i, j)$th entry is $M_d(y)[i, j] = l_y(z^{i+j}) = y_{i+j}$, $\forall i, j \in \mathbb{N}^p_2$.

Similarly, the localizing matrix of order $d$ associated with a sequence $(y_\alpha)_\alpha$ and a polynomial $g(z)$ is the matrix $M_d(g y)$ such that $l_y(g(z)p(z)^2) = p'M_d(g y)p$ for all polynomial $p(z)$ of degree $d$.

**Proposition 4.11.** Let $Z$ be a compact basic semi-algebraic set defined by $Z = \{z \in \mathbb{R}^p \mid g_k(z) \geq 0, k = 1, \ldots, n_Z\}$. Then, a necessary condition for a sequence $(y_\alpha)_\alpha$ to have a representing measure $\mu \in \mathcal{M}^+(z)$ is

$$M_d(y) \geq 0, \quad M_d(g_k y) \geq 0, \quad \forall d \in \mathbb{N}, \forall k = 1, \ldots, n_Z.$$

Finally, we introduce

$$[0, T] \times Q \times U = \{(t, q, u) \mid q = (q_{11}, q_{12}, q_{21}, q_{22}), g_1(t, q, u) := t(T - t) \geq 0,\]

$$g_2(t, q, u) := 1 - q_{11}^2 - q_{12}^2 \geq 0, \quad g_3(t, q, u) := 1 - q_{21}^2 - q_{22}^2 \geq 0,\]

$$g_4(t, q, u) := 1 - u^2 \geq 0\},$$

and

$$Q_f = \{(t, q) \in \mathbb{R}^5 \mid g_0^f(t) := t(T - t) \geq 0, g_1^f(q) := q_{11} = 0,\]

$$g_2^f(q) := q_{12} = 0, \quad g_3^f(q) := q_{21} = 0, \quad g_4^f(q) := q_{22} = 0\}.$$ 

We denote by $l_{y^{\mu}}$, $l_{y^{\mu\prime}}$ the Riesz functionals associated respectively with the sequences $y^{\mu}$ and $y^{\mu\prime}$. Then, the moment SDP problem associated with (4.1) is

$$T_{SDP} = \inf_{y^{\mu} \in y^{\mu\prime}, \mu \prime} l_{y^{\mu\prime}}(1)$$

$$l_{y^{\mu}} \left( \frac{\partial v}{\partial t} + \frac{\partial v}{\partial q} (F + uG) \right) = l_{y^{\mu\prime}}(v(\cdot, q_f)) - v(0, q_0), \quad \forall v \in \mathbb{R}[t, q],$$

$$M_d(y^{\mu}) \geq 0, \quad M_d(g_i y^{\mu}) \geq 0, \quad i = 1, \ldots, 4, \quad \forall d \in \mathbb{N},\]

$$M_d(y^{\mu\prime}) \geq 0, \quad M_d(g_i^f y^{\mu\prime}) \geq 0, \quad i = 0, \ldots, 4, \quad \forall d \in \mathbb{N}.$$ 

At the end, according to Proposition 4.11, we have $T_{LP} \geq T_{SDP}$ and this is in fact an equality according to Theorem 3.8 from [24].
Step 3: Hierarchy of SDP problems. Note that $M_{d+1}(y) \geq 0$ implies $M_d(y) \geq 0$. The LMI constraints and the sequence $(g_n)$ of (4.4) are truncated which lead to $r \geq 1$ to the Lasserre’s hierarchy parameterized by $r \geq 1$

$T_{LMI}^r = \inf_{(y^r)_{|\omega| \leq 2r} \in \mathcal{M}} \{ l_{y^r}(1) \}$

$l_{y^r} \left( \frac{\partial v}{\partial t} + \frac{\partial v}{\partial q} (F + uG) \right) = l_{y^r} (v(\cdot, q)) - v(0, q_0), \ \forall v \in \mathbb{R}[t, q]$. (4.5)

$M_r(y^r) \geq 0, M_{r-s_i}(g_i y^r) \geq 0, \ i = 1 \ldots 4$

$M_r(y^{\mu^r}) \geq 0, M_r(g_i y^{\mu^r}) \geq 0, \ i = 0 \ldots 4.$

where $s_i = \deg(g_i)/2$ if $\deg(g_i)$ is even and $s_i = (\deg(g_i) + 1)/2$ otherwise. The main result is then the following.

Proposition 4.12 (Theorem 5.6, [24]). We have

$T_{\min} = T_{LP} = T_{SDP} \geq \ldots \geq T_{LMI}^{r+1} \geq T_{LMI}^r \geq \ldots \geq T_{LMI}^1.$

Moreover the sequence of lower bounds $(T_{LMI}^r)$ converges to $T_{\min}$ as $r \to \infty$.

Summary of the LMI method. The moment/LMI method approach for optimization consist in reformulating an optimization problem as a linear program on measures. When the data is polynomial, a hierarchy of LMI relaxations can be constructed, whose costs converge to that of the original problem. The strong feature of the method is that those LMI generate lower bounds on the true cost, and can therefore be used as certificates of global optimality. On the other hand, the weak points of the method are its poor algorithm complexity for unstructured problem, as well as for the special case of optimal control, the unavailability of a generic method to recover controls. Note that the passage to a given LMI relaxation starting from measure problem (4.1) or (4.3) can be fully automated with high-level commands using the GloptiPoly toolbox [19].

4.3.3. Validation of the numerical results. The problem (4.5) corresponds to the multisaturation problem of two spins associated with the LP problem (4.1). Likewise, we can construct from the LP problem (4.3) a hierarchy of LMI relaxations parameterized by $r$ in the single spin case and the two spins case to compare the two formulations. We use the Mosek toolbox [29] to solve the SDP problems. Let $T_f$ denote the best solution found with the Bocop software, given in Table 2 for the single spin case and the two spins case. The value of the parameters for (4.5) are: $\varepsilon = 0.1, q_0 = (0, 1, 0, 1), q_f = (0, 0, 0, 0)$ and $T = t_f$. The lower bounds of $T_{\min}$ are $T_{LMI}^r$ (resp. $T_{LMI}^r$) associated with the LP problem (4.1) (resp. (4.3)) and are given in Table 3.

We denote by $n = \dim Q, m = \dim U$ and $n_d = 2$ is the number of modes for the second formulation. At the relaxation order $d$, the number of moments involved in the SDP problem associated with (4.1) and (4.3) is given by $N_m = \left( \frac{n+m+1+2d}{n+m+n+1} \right)$ and $N_m = \left( \frac{n+1+2d}{n+1} \right)$ respectively. In Fig. 16 are represented the relative error $\text{err}(r) = (t_f - T_{LMI})/t_f$ for the cases $C_1, C_2, C_3$ and $C_4$, where $T_{LMI}$ is the optimal value of (4.5) in the single spin case and the two spins case. Note that in the single spin case, we know the structure and the optimal value of the global solution and we can compute the numerical gap between the direct approach and the moment/LMI approach.

Both formulations are computationally demanding, the relative errors on the final
<table>
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<tr>
<th>r</th>
<th>$N_m$</th>
<th>$(t_f - T_{LMI})/t_f$</th>
<th>$N_m$</th>
<th>$(t_f - T'_{LMI})/t_f$</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
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<td>2040</td>
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</tr>
<tr>
<td>8</td>
<td>5814</td>
<td>0.0461</td>
<td>2907</td>
<td>0.0716</td>
</tr>
</tbody>
</table>

Table 3. Single spin saturation for the case $C_2$. $T'_{LMI}$ is the optimal value of the hierarchy (4.5). Likewise, $T_{LMI}$ is the optimal value of the hierarchy of SDP problems derived from (4.3). $t_f = 110.44$ is the best time found by the Bocop software. The second and fourth columns are the relative errors between the best solution found by the Bocop software and the one found by moment/LMI techniques for each relaxation order $r$.

Times found by Bocop for the cases $C_2$, $C_3$, $C_4$ are less than 5% for the single spin saturation and less than 10% for the multisaturation. Note that these two formulations have to be compared not only on the sharpness of the lower bounds but also considering the number of moments involved in the hierarchy.

Figure 16. Saturation problem of one spin and two spins for the cases $C_1$, $C_2$, $C_3$ and $C_4$. Relative error $err(r) = (t_f - T'_{LMI})/t_f$ where $r$ is the order of relaxation, $T'_{LMI}$ is the optimal value of (4.5) using the formulation (4.3) and $t_f$ is the final time computed with Bocop.
Influence of the parameters on the BC-extremals. We use a combination of multiple shooting and differential path following methods to analyze the influence of the parameters on BC-extremals from problem (P_{BS}). We recall the multiple shooting technique in section 4.4.1, we give some details about homotopy and monitoring in section 4.4.2 and we present the results in section 4.4.3.

4.4.1. Multiple shooting (HamPath). Since the optimal structures are composed of sequences of bang and singular arcs, we must use multiple shooting instead of single shooting. We refer to [14, 28] for details about multiple shooting algorithms and to [8] and [9] for explanations about multiple shooting in the context of medical imaging. One particularity is that the solutions end with a singular arc and not a bang arc, contrary to the examples given in [28]. Because of that, the shooting equations are more intricate as we can see it hereinafter.

Let \( u_s(x, p) \) denote the singular control and \( u_\pm = \pm u_{\text{max}} = \pm 1 \) the positive and negative bang controls. Let us assume we have a solution with a structure of the form BSBS, i.e. Bang-Singular-Bang-Singular. We note \( y = (p_0, t_f, t_1, t_2, t_3, z^1, z^2, z^3) \) the unknowns of the shooting function which is given by:

\[
S: \mathbb{R}^{32} \rightarrow \mathbb{R}^{32}
\]

\[
y = \begin{bmatrix} p_0 \\ t_f \\ t_1 \\ t_2 \\ t_3 \\ z^1 \\ z^2 \\ z^3 \end{bmatrix} \quad \mapsto \quad S(y) = \begin{bmatrix} u_\pm H_1(z^0) + p^0 \\ H_1(z^1) \\ H_{01}(z^1) \\ H_1(z^3) \\ H_{01}(z^3) \\ y_2(t_f, t_3, z^3, u_s) \\ z_2(t_f, t_3, z^3, u_s) \\ (p_z_1(t_f, t_3, z^3, u_s) + p_{zz}(t_f, t_3, z^3, u_s)) \gamma + p^0 \\ z(t_1, 0, z^0, u_\pm) - z^1 \\ z(t_2, t_1, z^1, u_s) - z^2 \\ z(t_3, t_2, z^2, u_\pm) - z^3 \end{bmatrix}
\]

where \( p^0 = -1 \) in the normal case, where \( z^0 = (q_0, p_0) \) is the initial state-costate vector with \( q_0 = (0, 1, 0, 1) \). To get a BC-extremal we want to solve the shooting equations

\[
S(y) = 0.
\]

The first equation comes from the fact that the final time is free, the four following equations means that the associated extremal becomes singular at \( z^1 \) and \( z^3 \). The last three matching equations improve numerical stability. Note that we have only three equations (the three remaining equations) associated to the final condition: \( q(t_f) = (0, 0, 0, 0) \). This is because we can find a redundant equation and this is due to the fact that the trajectory ends with a singular (and not bang) arc. Indeed, the fact that \( H_1(z^3) = H_{01}(z^3) = 0 \) implies that \( H_1(z(t_f, t_3, z^3, u_s)) = 0 \)
Hence, for any zero of the shooting function, the associated trajectory reaches the target $q_f = (0,0,0,0)$ at the final time if $(p_{y_1}(t_f,t_3,z^3,u_3), p_{z_1}(t_f,t_3,z^3,u_3)) \neq 0$. Let us recall that one difficulty to solve a shooting equation is to have a good initial guess. To determine the structure and make the shooting method converge, we use direct methods from the Bocop software. This combination of direct and indirect methods has already been successful in the contrast problem [8] by nuclear magnetic resonance in medical imaging.

4.4.2. Homotopy and monitoring. Once we have a solution obtained by a multiple shooting method, we can use differential homotopy techniques [1] to study the deformation of the solution with respect to the relaxation parameters. The homotopy method, from HamPath software [15], is based on Predictor-Corrector algorithm with a high order and step-size control Runge-Kutta scheme for the prediction and with a classical simplified Newton method for the correction. We combine the differential path following method with monitoring at each accepted step of the integration to detect if structural changes occur during the homotopy. We consider three different monitoring for which we give the associated action for the saturation problem:

- check if each arc (except the first and the last arcs) has positive length, that is if $t_i \leq t_{i+1}$. If not, the arc with negative length has to be removed.
- check if the singular control on each singular arc (except the last arc because it is 0) is admissible. If not, one bang arc has to be added.
- check if the switching function on each bang arc remains of constant sign.

Let us illustrate the third monitoring on an example. We write $\Gamma = \rho \cos \theta$ and $\gamma = \rho \sin \theta$, and we consider an homotopy (called H1a) on $\theta$ with $\rho = \tilde{\rho} \approx 0.0551$, from $\theta = \theta_{\text{max}} = \text{atan}(2) \approx 1.1071$ to $\theta = \theta_{\text{min}} = 0.02$. One starts from $\theta = \theta_{\text{max}}$ with a structure of the form BSBS ($\sigma_- \sigma_+ \sigma_0 \sigma_0$). Around $\theta = \theta_{1a,1} \approx 0.5069$, the monitoring detects a change in the structure. Let us denote by $\theta_{1a,1}^+ \geq \theta_{1a,1} \geq \theta_{1a,1}^-$ the two values of $\theta$ at the two consecutive steps such that these inequalities are satisfied. Since, for $\theta = \theta_{1a,1}^-$, the switching function crosses 0 two times on the first bang arc, one has to stop the homotopy and add a singular arc, see figure 17.

4.4.3. Numerical results. We write $\Gamma = \rho \cos \theta$ and $\gamma = \rho \sin \theta$, and we consider only homotopies on $\theta$ with $\rho = \tilde{\rho} \approx 0.0551$. This particular value of $\rho$ is such that we retrieve the fat case, that is $T_1 = 0.2$ and $T_2 = 0.1$, with $h_{\text{max}} = 1$ and $\omega_{\text{max}} = 2\pi \times 32.3$, this specific value of $\omega_{\text{max}}$ being excerpted from [11]. We present some results for a range of values of $\theta$ to illustrate the role of this parameter on the structure of some BC-extremals from problem $P_{\text{BS}}$. The value of $\theta$ ranges between $\theta = \theta_{\text{max}} = \text{atan}(2) \approx 1.1071$ and $\theta = \theta_{\text{min}} = 0.02$. The maximal bound is chosen to satisfy the physical constraint $0 < \gamma \leq 2\Gamma$ while the minimal bound is chosen
to include all the practical cases, see Table 1. The saturation of a pair of spins is a much more complex problem than the mono-saturation, so we do not intend to get any optimal synthesis but just optimal trajectories to steer both spins from the North Pole to the center of the Bloch ball in minimum time and in a synchronized fashion. Besides, we just analyze the influence of $\theta$ for fixed values of $\rho = \bar{\rho}$ and $\varepsilon = 0.1$.

The methodology is the following. For a given initial value of $\theta$, we use direct method to determine the structure and to initialize the multiple shooting method. Then, we perform homotopies on $\theta$ with monitoring to stop it if necessary. If a change in the structure is detected, then, we update the multiple shooting function and continue the homotopy until we reach the final value of $\theta$ if possible. Note that in the multi-saturation problem, there exist many local solutions. Thus, for a given value of $\theta$ we must compare the cost of each solution. This leads to compute several path of zeros and then compare them in terms of cost. We choose to present four different paths, see Table 4 and Fig 21. Note that the paths of zeros are computed with a very good accuracy according to the bottom-right subgraph of Fig 21. Before we give some details about the homotopies, we present three particular cases excerpted from these homotopies: the Water case ($\omega_{\text{max}} = 10.2684$), the Fat case ($\omega_{\text{max}} = 202.9469 \approx 2\pi \times 32.3$) and the Cerebrospinal Fluid case ($\omega_{\text{max}} = 61.1840$). See Table 1 for the corresponding values of $\theta$. All the others cases have the same optimal structure than the Cerebrospinal Fluid case. One can see from Fig 21 that for the Water case, the optimal structure is of the form $\sigma \cdots \sigma \sigma_0$ and the solution is given by the homotopy (H1a). The solution is given on Fig. 18. For the Fat case, the structure is of the form $\sigma \cdots \sigma \sigma_0 \sigma_0$ and it is also given by (H1a), see Fig. 19. Finally, for the Cerebrospinal Fluid case, the structure is of the form $\sigma \cdots \sigma_0 \sigma_0 \sigma_0 \sigma_0$ and it is given by (H1b), see Fig. 20. One can notice different interactions between bang and singular arcs depending on the
case, and one can see that the horizontal line $z = z_s = \gamma/2\delta$ plays a crucial role in the optimal trajectory when it intersects the Bloch ball.

Figure 18. Water case: $\rho = \overline{\rho}$, $\theta = 0.7854$ and $\varepsilon = 0.1$. Trajectories of spins 1 and 2, control and $H_1$.

Figure 19. Fat case: $\rho = \overline{\rho}$, $\theta = 0.4636$ and $\varepsilon = 0.1$. Trajectories of spins 1 and 2, control and $H_1$. 
Figure 20. Fluid: \( \rho = \bar{\rho}, \theta = 0.1489 \) and \( \varepsilon = 0.1 \). Trajectories of spins 1 and 2, control and \( H_1 \).

Let us give now some details about the different homotopies. The first two paths (called \( H_{1a} \) and \( H_{1b} \)) are interesting since, even if they are distinct paths (see bottom-left subgraph of Fig 21), they intersect in terms of cost for a value \( \theta^* \in [\theta_{1b,1}, \theta_{1b,2}] \), where \( \theta_{1b,1} \approx 0.2752 \) and \( \theta_{1b,2} \approx 0.3018 \). For \( \theta \leq \theta^* \), \( H_{1b} \) is better and for \( \theta \geq \theta^* \), \( H_{1a} \) is better, see top-left subgraph of Fig 21. The homotopy \( H_{1a} \) has been already introduced in section 4.4.2. One starts from \( \theta = \theta_{\text{max}} \) with a structure of the form \( \text{BSBS} (\sigma^- \sigma^s, \sigma^+, \sigma_0) \). Around \( \theta = \theta_{1a,1} \approx 0.5069 \), the monitoring detects a change in the structure: a singular arc has to be added inside the first negative bang arc and the structure becomes \( \sigma^- \sigma^s \sigma^- \sigma^+ \sigma_0 \). A new change occurs around \( \theta_{1a,2} \approx 0.2722 \): the second singular arc vanishes and the structure becomes \( \sigma^- \sigma^s \sigma^+ \sigma_0 \). We stop the homotopy here since the homotopy \( H_{1b} \) is better for this value of \( \theta \). Let us explain the homotopy \( H_{1b} \): this homotopy starts from \( \theta = \theta_{\text{min}} \) with a structure of the form \( \text{BSBSBS} (\sigma^- \sigma^s, \sigma^+ \sigma_0) \). A first change in the structure is detected around \( \theta_{1b,1} \approx 0.2752 \): the second singular arc vanishes and the structure becomes \( \sigma^- \sigma^s \sigma^+ \sigma_0 \). A second change occurs around \( \theta_{1b,2} \approx 0.3018 \): the control saturates at the end of the first singular arc. We denote by \( \sigma^-_s \) this singular arc with since at the end the control takes the value \( -1 \). The structure is now \( \sigma^- s \sigma^+ \sigma_0 \) and we do not continue the homotopy since \( H_{1a} \) is better for \( \theta = \theta_{1b,2} \). The different structures with the names of the homotopies and the associated figures to observe the trajectories and the control are given in Table 4 and Figs. 22–36.

Let us give some details about the two last homotopies. The homotopy \( H_2 \) starts with a local solution (not globally optimal, see top-right subgraph of Fig 21) of the form \( \sigma^+ \sigma^s \sigma^+ \sigma^+ \sigma_0 \) for \( \theta = 0.2 \). This solution has the particularity that all the bang arcs are positive bang arcs. Besides, the trajectory (see Fig 29) has a self-intersection which prevents the BC-extremal to be globally optimal. During
the homotopy there is only one change, the first singular arc vanishes around $\theta = \theta_{2,1} \approx 0.3618$. This homotopy may be compared with the homotopy H3. The homotopy H3 is similar as H2 but the last bang arc is longer. During this last bang arc, the trajectory realizes a complete turn around its center. There is also one single change around $\theta = \theta_{2,1} \approx 0.4778$, from which the first singular arc vanishes.

<table>
<thead>
<tr>
<th>Name</th>
<th>Init</th>
<th>Transition</th>
<th>End</th>
</tr>
</thead>
<tbody>
<tr>
<td>H1a</td>
<td>$\theta_{\text{max}} \approx 1.1071$</td>
<td>$\theta_{1a,1} \approx 0.5069$</td>
<td>$\theta_{1a,2} \approx 0.2722 &lt; \theta_{\text{min}} = 0.02$</td>
</tr>
<tr>
<td></td>
<td>$\sigma_– \sigma_+ \sigma_0$</td>
<td>$\sigma_– \sigma_+ \sigma_0$</td>
<td>$\sigma_– \sigma_+ \sigma_0$</td>
</tr>
<tr>
<td></td>
<td>fig 22</td>
<td>fig 23</td>
<td>fig 24</td>
</tr>
<tr>
<td>H1b</td>
<td>$\theta_{\text{min}}$</td>
<td>$\theta_{1b,1} \approx 0.2752$</td>
<td>$\theta_{1b,2} \approx 0.3018 &lt; \theta_{\text{max}}$</td>
</tr>
<tr>
<td></td>
<td>$\sigma_– \sigma_+ \sigma_0$</td>
<td>$\sigma_– \sigma_+ \sigma_0$</td>
<td>$\sigma_– \sigma_+ \sigma_0$</td>
</tr>
<tr>
<td></td>
<td>fig 25</td>
<td>figs 26 and 27</td>
<td>fig 28</td>
</tr>
<tr>
<td>H2</td>
<td>$\theta = 0.2$</td>
<td>$\theta_{2,1} \approx 0.3618$</td>
<td>$\theta = \text{atan}(1.5) \approx 0.9828$</td>
</tr>
<tr>
<td></td>
<td>$\sigma_+ \sigma_+ \sigma_0$</td>
<td>$\sigma_+ \sigma_+ \sigma_0$</td>
<td>$\sigma_+ \sigma_+ \sigma_0$</td>
</tr>
<tr>
<td></td>
<td>fig 29</td>
<td>figs 30 and 31</td>
<td>fig 32</td>
</tr>
<tr>
<td>H3</td>
<td>$\theta = 0.25$</td>
<td>$\theta_{3,1} \approx 0.4778$</td>
<td>$\theta = \theta_{\text{max}}$</td>
</tr>
<tr>
<td></td>
<td>$\sigma_+ \sigma_+ \sigma_0$</td>
<td>$\sigma_+ \sigma_+ \sigma_0$</td>
<td>$\sigma_+ \sigma_+ \sigma_0$</td>
</tr>
<tr>
<td></td>
<td>fig 33</td>
<td>figs 34 and 35</td>
<td>fig 36</td>
</tr>
</tbody>
</table>

**Table 4.** The homotopies are detailed on each line. The first column gives the name of the homotopy. The second column “Init” gives the initial value of $\theta$, the associated structure with the reference to the figure which presents the trajectory with the control and the switching function. For this initial value, the solution is obtained from direct method and multiple shooting. The column “Transition” gives the same details but when a first change in the structure is detected during the homotopy thanks to the monitoring. The last column “End” gives again the same details at the end of the homotopy. For the homotopies H1a and H1b, the end occurs when a second change in the structure is detected while for H2 and H3, the end occurs when the final value of $\theta$ is reached. The red and green colors may help to understand the change in the structure, which is explained more in details in this section.
Figure 21. The homotopies $H_{1a}$, $H_{1b}$, $H_2$ and $H_3$ are presented respectively in blue, red, black and black. The homotopies $H_2$ and $H_3$ are presented only on the top-right subgraph. The plain and dashed lines distinguish the different structures. The top-left subgraph gives the cost (i.e. the final time $t_f$) with respect to $\theta$ for the homotopies $H_{1a}$ and $H_{1b}$. One can see (it is more visible in the zoom) the intersection of the two paths of zeros in terms of cost. $H_{1b}$ is better for small value of $\theta$ and $H_{1a}$ is better for greater values. One can notice that this two paths of zeros are distinct from the bottom-left subgraph. This subgraph gives the norm of the initial adjoint vector with respect to the homotopic parameter. The norm of the shooting function along the paths $H_{1a}$ and $H_{1b}$ is given in the bottom-left subgraph. One can see the very good accuracy. The strategies from $H_{1a}$ and $H_{1b}$ are compared with homotopies $H_2$ and $H_3$ on the top-right subgraph.
Figure 22. H1a: $\rho = \bar{\rho}, \theta = \theta_{\text{max}}$ and $\varepsilon = 0.1$. Trajectories of spins 1 and 2, control and $H_1$.

Figure 23. H1a: $\rho = \bar{\rho}, \theta = \theta_{10,1}$ and $\varepsilon = 0.1$. Trajectories of spins 1 and 2, control and $H_1$.

Figure 24. H1a: $\rho = \bar{\rho}, \theta = \theta_{10,2}^+$ and $\varepsilon = 0.1$. Trajectories of spins 1 and 2, control and $H_1$. 
Figure 25. H1b: \( \rho = \bar{\rho}, \theta = \theta_{\min} \) and \( \varepsilon = 0.1 \). Trajectories of spins 1 and 2, control and \( H_1 \).

Figure 26. H1b: \( \rho = \bar{\rho}, \theta = \theta_{1b,1}^- \) and \( \varepsilon = 0.1 \). Trajectories of spins 1 and 2, control and \( H_1 \).

Figure 27. H1b: \( \rho = \bar{\rho}, \theta = \theta_{1b,1}^+ \) and \( \varepsilon = 0.1 \). Trajectories of spins 1 and 2, control and \( H_1 \).

Figure 28. H1b: \( \rho = \bar{\rho}, \theta = \theta_{1b,2}^+ \) and \( \varepsilon = 0.1 \). Trajectories of spins 1 and 2, control and \( H_1 \).
Figure 29. H2: $\rho = \bar{\rho}$, $\theta = 0.2$ and $\varepsilon = 0.1$. Trajectories of spins 1 and 2, control and $H_1$.

Figure 30. H2: $\rho = \bar{\rho}$, $\theta = \theta_{2,1}$ and $\varepsilon = 0.1$. Trajectories of spins 1 and 2, control and $H_1$.

Figure 31. H2: $\rho = \bar{\rho}$, $\theta = \theta_{2,1}^{-1}$ and $\varepsilon = 0.1$. Trajectories of spins 1 and 2, control and $H_1$.

Figure 32. H2: $\rho = \bar{\rho}$, $\theta = \text{atan}(1.5)$ and $\varepsilon = 0.1$. Trajectories of spins 1 and 2, control and $H_1$. 
Figure 33. $H_3$: $\rho = \bar{\rho}$, $\theta = 0.25$ and $\varepsilon = 0.1$. Trajectories of spins 1 and 2, control and $H_1$.

Figure 34. $H_3$: $\rho = \bar{\rho}$, $\theta = \theta_3^{-1}$ and $\varepsilon = 0.1$. Trajectories of spins 1 and 2, control and $H_1$.

Figure 35. $H_3$: $\rho = \bar{\rho}$, $\theta = \theta_3^1$ and $\varepsilon = 0.1$. Trajectories of spins 1 and 2, control and $H_1$.

Figure 36. $H_3$: $\rho = \bar{\rho}$, $\theta = \text{atan}(2)$ and $\varepsilon = 0.1$. Trajectories of spins 1 and 2, control and $H_1$. 
5. Algebraic computations for multisaturation with $B_1$-inhomogeneity

The Maple symbolic software is used to perform algebraic computations related to singularity analysis of the extremal trajectories, in particular in relation with the determination of bridges. The computations boil down to the computation of Gröbner bases and one needs the following concepts and techniques of this area.

5.1. Operation on polynomial ideals. Let $n$ be a positive integer and we consider polynomials in $A = \mathbb{C}[X_1, \ldots, X_n]$. Given a system of equations $\mathcal{F}$, one considers the ideal $I$ coding the set $V(\mathcal{F})$ of zeros of $\mathcal{F}$.

Given an ideal $J$ and a polynomial $F$, saturating $I$ by $F$ means computing a set of generators of the ideal $$(I : F^\infty) = \{ g \in A, \exists m \in \mathbb{N}, g F^m \in I \}.$$ The radical of an ideal $I$ is the set: $$\sqrt{I} = \{ f \in A, \exists m \in \mathbb{N}, f^m \in I \}$$ and it has the same set of zeros. An operation used frequently is, given a set of generators of $I$, computing a set of generators of $J \supset I$ such that $\sqrt{I} = \sqrt{J}$. This can be done by computing the square free form of the generators: given a polynomial $f$ with decomposition in primes $p_1^{n_1} \cdots p_r^{n_r}$, the square free form is $\text{sqfr}(f) = p_1 \cdots p_r$.

5.2. Frame curves. Frame curves associated to the saturation of a single spin lead to the following.

- **Collinearity locus:** $C$ is defined as the set where $f$ and $g$ are linearly dependent. Outside zero, it is defined by: $\exists \lambda$ such that $F = \lambda G$, that is:

$$- \Gamma y_1 = \lambda z_1, \quad \gamma(1 - z_1) = \lambda y_1,$$
$$- \Gamma y_2 = \lambda z_2 (1 - \varepsilon), \quad \gamma(1 - z_2) = \lambda y_2 (1 - \varepsilon),$$

The projections on $q_i$-spaces are the ovals:

$$\Gamma_i y^i_1 = \gamma (1 - z_i) z_i, \quad 0 \leq z_i \leq 1, \quad i = 1, 2$$

interacted with one of the sets

$$\Gamma(1 - \varepsilon) y_1 z_2 = \Gamma y_2 z_1, \quad (1 - \varepsilon) y_2 (1 - z_1) = y_1 (1 - z_2)$$

- **Singularity locus:** $S$ is defined as the set where $G$ and $[F, G]$ are linearly independent. Outside zero, it is defined by: $\exists \lambda$ such that $[F, G] = \lambda G$, that is

$$\delta z_1 - \gamma = -\lambda z_1, \quad \delta y_1 = \lambda y_1,$$
$$\delta z_2 - \gamma = -\lambda z_2, \quad \delta y_2 = \lambda y_2,$$

Projections on each $q_i$-spaces will form the two singular lines

$$z_i = \gamma / 2 \delta, \quad y_i = 0, \quad i = 1, 2.$$ The additional relations define the full locus.

5.3. Singularity classification: exceptional case $H_F = 0$. 


Conventions and notations. For the computations, we use the translation
\[ z_1 \leftrightarrow z_1 + 1, \quad z_2 \leftrightarrow z_2 + 1, \]
which places the center of the coordinates at the North pole of the Bloch ball. In this new system of coordinates, the center of the Bloch ball has coordinates \((0, -1, 0, -1)\).

We have in the exceptional case the constraint
\[ p \cdot F = p \cdot G = p \cdot [F, G] = 0, \]
hence \(p\) can be eliminated in the relation defining the control
\[ p \cdot ([[G, F], F] + u [[G, F], G]) = 0 \]
which is computed as the feedback
\[ u_s(q) = -\frac{D'(q)}{D(q)} \]
with
\[ D = \det(F, G, [[G, F], [[G, F], G]]) \]
\[ = \det\begin{bmatrix} -\Gamma y & -z - 1 & \delta z - \Gamma & 2\delta y \\ -\gamma z & y & \delta y & -2\delta z + \Gamma - \delta \\ -\Gamma y (1 - \epsilon)(-z - 1) & (1 - \epsilon)(\delta z - \Gamma) & (1 - \epsilon)^2(2\delta y) \\ -\gamma z y (1 - \epsilon) & (1 - \epsilon)\delta y & (1 - \epsilon)^2(-2\delta z + \Gamma - \delta) \end{bmatrix} \]
and
\[ D' = \det(F, G, [G, F], [[G, F], F]) \]
\[ = \det\begin{bmatrix} -\Gamma y & -z - 1 & \delta z - \Gamma & \gamma(\gamma - 2\Gamma) + \delta^2(z + 1) \\ -\gamma z & y & \delta y & \delta^2 y \\ -\Gamma y (1 - \epsilon)(-z - 1) & (1 - \epsilon)(\delta z - \Gamma) & (1 - \epsilon)^2(\gamma(\gamma - 2\Gamma) + \delta^2(z + 1)) \\ -\gamma z y (1 - \epsilon) & (1 - \epsilon)\delta y & (1 - \epsilon)^2\delta^2 y \end{bmatrix} \]

Using a time reparameterization, this leads to analyze the \(C^\omega\)-vector field in the \(q\)-space
\[ X = DF - D'G. \]

The following polynomials will appear frequently in the remainder of the section:
- \( P_{y_1} := y_1 - (1 - \epsilon)y_2 \)
- \( P_{y_2} := y_2 - (1 - \epsilon)y_1 \)
- \( P_{z_1} := 2(\Gamma - \gamma)z_1 + 2\Gamma - \gamma \)
- \( P_{z_2} := 2(\Gamma - \gamma)z_2 + 2\Gamma - \gamma \)

The root of the univariate polynomials \( P_{z_1} \) and \( P_{z_2} \), in \( z_1 \) and \( z_2 \) respectively, is
\[ z_s = \frac{\gamma - 2\Gamma}{2\Gamma - 2\gamma}. \]
Figure 37. Structure of the study of the singularities of \( \{ D = 0 \} \)

Singularities of \( \{ D = 0 \} \).

**Proposition 5.1.** The set of points satisfying \( D = \frac{\partial D}{\partial y_1} = \frac{\partial D}{\partial y_2} = \frac{\partial D}{\partial z_1} = \frac{\partial D}{\partial z_2} = 0 \) is given, generically on authorized values of \( \gamma, \Gamma \), by

1. the point \( y_1 = y_2 = z_1 = z_2 = 0 \), and
2. the curve defined by \( P_{y_1} = P_{z_1} = P_{z_2} = 0 \), which is parameterized by \( y_2 \) as

\[
\begin{cases}
y_1 = (1 - \epsilon)y_2 \\
z_1 = z_2 = z_S = \frac{\gamma - 2\Gamma}{2\Gamma - \gamma}.
\end{cases}
\]

If \( \gamma = \Gamma \) (for example if the matter is water), only the former solution exists.

**Proof.** The structure of this proof is summarized in Fig. 37. The determinant \( D \) can be factored as \( (1 - \epsilon)\tilde{D} \). The singularities of \( D \) and those of \( \tilde{D} \) are the same, so for the study, we consider the ideal

\[
I := \left\langle \frac{\partial \tilde{D}}{\partial y_1}, \frac{\partial \tilde{D}}{\partial y_2}, \frac{\partial \tilde{D}}{\partial z_1}, \frac{\partial \tilde{D}}{\partial z_2} \right\rangle.
\]

In order to eliminate \( y_1, y_2 \) and \( z_1 \) from the ideal \( I \), we compute a Gröbner basis \( G \) of \( I \) with respect to the elimination ordering \( y_1 > y_2 > z_1 > z_2 > \epsilon > \Gamma > \gamma \). This computation yields that

\[
I \cap \mathbb{Q}[z_2, \epsilon, \Gamma, \gamma] = \left\langle (\epsilon^2 - 2)\tilde{z}^2(2\Gamma - \gamma)(\Gamma - \gamma)z_2^3P_{z_2}^3 \right\rangle,
\]

so singular points necessarily satisfy

\[
\begin{cases}
z_2 = 0 \\
\text{or} \\
P_{z_2} = 0 \iff z_2 = \frac{\gamma - 2\Gamma}{2\Gamma - \gamma}.
\end{cases}
\]
If $\Gamma = \gamma$ (that is, if the matter is water), the second of these solutions does not exist. If $\gamma = 2\Gamma$ (which means that the matter is on the limit of the domain of validity $2\Gamma \geq \gamma$), both solutions coincide.

In all other cases, there are 2 distinct possible values for $z_2$, and we consider both cases: we consider the two ideals

$$I_1 := \langle \text{sqfr}(G), z_2 \rangle$$
$$I_2 := \langle \text{sqfr}(G), P_{z_2} \rangle$$

where for any polynomial $f$, $\text{sqfr}(f)$ is the square-free part of $f$ and $\text{sqfr}(G)$ means that we apply $\text{sqfr}$ to each element of $G$.

In order to lift the partial solution $z_2 = 0$, we compute a Gröbner basis $G_1$ of $I_1$ with respect to the ordering $y_1 > y_2 \gg z_1 > z_2 > \epsilon > \Gamma > \gamma$, and we find that this ideal contains $\gamma z_1^2(\epsilon - 1)^2(2\Gamma - \gamma)$, so $z_1 = 0$.

We then compute a Gröbner basis of $\langle \text{sqfr}(G_1), z_1 \rangle$ with respect to the order $y_1 \gg y_2 > z_1 > z_2 > \epsilon > \Gamma > \gamma$, and we find that this ideal contains $\Gamma \gamma y_2^2(\Gamma - \gamma)(2\Gamma - \gamma)^2(\epsilon - 2)$, so $y_2 = 0$.

Finally, adding $y_2$ to the ideal yields that $0 = \Gamma y_1(\epsilon - 1)(2\Gamma - \gamma)$, so the complete solution is

$$(y_1, y_2, z_1, z_2) = (0, 0, 0, 0).$$

We now consider the partial solution $z_2 = (\gamma - 2\Gamma)/(2\Gamma - 2\gamma)$. We compute a Gröbner basis $G_2$ of $I_2$ with respect to the order $y_1 \gg y_2 > z_1 > z_2 > \epsilon > \Gamma > \gamma$, and we find that the ideal contains $z_2 \gamma(z_1 - z_2)^2$.

Since this case was already studied, we may assume that $z_2 \neq 0$, so

$z_1 = z_2 = \frac{\gamma - 2\Gamma}{2\Gamma - 2\gamma}$.

Adding $z_1 - z_2$ to $\text{sqfr}(G_2)$ and computing a Gröbner basis for the order $y_1 \gg y_2 > z_1 > z_2 > \epsilon > \Gamma > \gamma$, we find that the ideal contains $\gamma^2 y_2(\epsilon - 1)P_{y_1}(2\Gamma - \gamma)$, so we have 2 new branches to consider.

If $y_2 \neq 0$, $y_1 = (1 - \epsilon)y_2$. Otherwise, by adding $y_2 = 0$ to the system of equations, we find that the ideal contains $\gamma y_1^2 z_2$, so $y_1 = 0$, and in particular, this point is on $\{P_{y_1} = 0\}$. □
\[ D = D' = 0 \]
\[ \hat{D} = 0 \]
\[ z_1 - z_2 = 0 \]
\[ D' = 0 \]
\[ (\epsilon - 1)y_2z_1 + y_1z_2 = 0 \]
\[ P_{y_2} = 0 \]
\[ p_2 = 0 \]
\[ y_1 = z_1 = 0 \]
\[ y_1 = 0 \]
\[ z_1 \neq 0 \]
\[ y_1 \neq 0 \]
\[ y_1 = y_2 = 0 \]
\[ a_3 \neq 0 \]
\[ y_2 = z_2 = 0 \]
\[ p_3 = 0 \]
\[ p_4 = 0 \]
\[ p_5 = 0 \]
\[ (\text{sol. 1}) \]
\[ (\text{sol. 2}) \]
\[ (\subset \text{ sol. 2}) \]
\[ (\text{sol. 3}) \]
\[ (\subset \text{ sol. 2}) \]
\[ (\text{sol. 4}) \]
\[ (\subset \text{ sol. 4}) \]
\[ (\subset \text{ sol. 2}) \]

Figure 38. Structure of the study of \( \{D = D' = 0\} \)

**Locus of \( \{D = D' = 0\} \).**

**Proposition 5.2.** The points of \( \{D = D' = 0\} \) are given by:

1. the plane
   \[ z_1 = z_2 = z_S = \frac{\gamma - 2\Gamma}{2\Gamma - 2\gamma} \] (5.5)

2. the line
   \[ \begin{cases} y_1 = y_2 = 0 \\ z_1 = z_2 \end{cases} \] (5.6)

3. the surface (parameterized by \( y_1, y_2 \))
   \[ z_1 = z_2 = \frac{\Gamma P_{y_2}^2 (\gamma - 2\Gamma)}{2(\Gamma - \gamma)a_3} \] (5.7)
   with
   \[ a_3 = (\Gamma + \gamma)P_{y_1}^2 + \epsilon(\epsilon - 2)\Gamma(y_1 - y_2)(y_1 + y_2) \]

4. the surface (parameterized by \( y_1, z_2 \))
   \[ y_2 = \frac{y_1 z_2}{(1 - \epsilon)z_1}, \quad z_1 = \frac{(2\Gamma - \gamma)z_2}{a_4} \] (5.8)
   with
   \[ a_4 = 2(\epsilon - 2)(\Gamma - \gamma)\epsilon z_2 + (2\Gamma - \gamma)(\epsilon - 1)^2 \]

5. the surface (parameterized by \( y_2, z_2 \))
   \[ z_1 = \frac{z_2 y_1}{(1 - \epsilon)y_2}, \quad y_1 = \frac{(1 - \epsilon)y_2 ((2\Gamma - \gamma)\Gamma y_2^2 + \gamma^2 z_2^2)}{a_5} \] (5.9)
   with
   \[ a_5 = \Gamma (2\epsilon(\epsilon - 2)(\Gamma - \gamma)z_2 + (\epsilon - 1)^2(2\Gamma - \gamma)) y_2^2 + \gamma^2 z_2^2 \]

**Proof.** The structure of this proof is summarized in Fig. 38. The determinant \( D' \) factors as
\[ D' = 2\gamma^2(2\Gamma - \gamma)(\Gamma - \gamma)(z_1 - z_2)(\epsilon - 1)((\epsilon - 1)y_2z_1 + y_1z_2). \]
so we form the two ideals

\[ I_1 = \langle \tilde{D}, z_1 - z_2 \rangle \]
\[ I_2 = \langle \tilde{D}, (\epsilon - 1)y_2 z_1 + y_1 z_2 \rangle . \]

If \( z_1 = z_2 \), after substitution, \( \tilde{D} \) has two factors depending on \( y_1, y_2, z_2 \): \( P_{z_2} \) and

\[ p_2 := 2(\Gamma - \gamma) ((\Gamma + \gamma) P_{y_1}^2 + \epsilon(\epsilon - 2)\Gamma(y_1 - y_2)(y_1 + y_2)) z_2 + \Gamma P_{y_2}^2 (2\Gamma - \gamma) \]

The polynomial \( P_{z_2} \) gives solution 1. Let

\[ a_3(y_1, y_2) = (\Gamma + \gamma) P_{y_1}^2 + \epsilon(\epsilon - 2)\Gamma(y_1 - y_2)(y_1 + y_2) \]

so that the coefficient of \( z_2 \) in \( p_2 \) is \( 2(\Gamma - \gamma)a_3 \), it is homogeneous in \( y_1, y_2 \) with degree 2. Its discriminant in \( y_2 \) is

\[-4(\epsilon - 2)^2 \epsilon^2 y_1^2 \gamma \Gamma . \]

Since the parameters \( \gamma, \Gamma \) are necessarily positive, this discriminant is negative, and thus the only real root of \( a_3(y_1, y_2) \) is \( y_1 = y_2 = 0 \). If \( y_1 = y_2 = 0 \), \( p_2 \) vanishes regardless of \( z_2 \).

If \( y_1 \neq 0 \), \( a_3(y_1, y_2) \) does not have any real root in \( y_2 \), and \( z_2 \) is given by

\[(z_1 =) z_2 = \frac{\Gamma P_{y_2}^2 (\gamma - 2\Gamma)}{2(\Gamma - \gamma)a_3(y_1, y_2)} . \]

We now turn to the other branch, defined by \((\epsilon - 1)y_2 z_1 + y_1 z_2 = 0\). If \( y_1 = z_1 = 0 \), there are 2 curves of singular points defined (in \( y_2, z_2 \)) by

\[ \Gamma(2\Gamma - \gamma)y_2^2 + \gamma^2 z_2^2 = 0 . \]

Since \( 2\Gamma \geq \gamma \), the only solution is \( z_2 = 0 \) with either \( 2\Gamma = \gamma \) (excluded) or \( y_2 = 0 \).

If \( y_1 = 0 \) and \( z_1 \neq 0 \), then (since \( \epsilon \neq 1 \)) we must have \( y_2 = 0 \). Furthermore, we may assume that \( z_1 \neq z_2 \) since this case was already studied. The remaining solutions form a curve defined by

\[ 0 = p_3 := (2(\epsilon - 2)(\Gamma - \gamma)ecz_2 + (2\Gamma - \gamma)(\epsilon - 1)^2)(z_1 - (2\Gamma - \gamma)z_2) . \]

Let \( a_4(z_1, z_2) \) be the coefficient of \( z_1 \) in \( p_3 \), the solutions are given by either

\[
\begin{align*}
    & a_4(z_2) \neq 0 \\
    & z_1 = \frac{2\Gamma - \gamma}{c_3(z_2)} z_2 
\end{align*}
\]

or (since by assumption \( z_1 \neq 0 \))

\[ z_2 = 2\Gamma - \gamma = 0 . \]

So we may assume that \( y_1 \neq 0 \). We compute a Gröbner basis of \( I_2 + \langle uy_1 - 1 \rangle \) for the order \( u \gg z_1 \gg y_1 \gg z_2 \gg y_2 \gg \epsilon \gg \gamma \gg \Gamma \). This basis contains a polynomial which factors as the product of

\[ p_4 = (2(\epsilon - 2)(\Gamma - \gamma)ecz_2 + (\epsilon - 1)^2 (2\Gamma - \gamma)) y_1 + (\epsilon - 1)(2\Gamma - \gamma)y_2 . \]
and
\[
p_5 = (\Gamma (2\epsilon (\epsilon - 2)(\Gamma - \gamma)z_2 + (\epsilon - 1)^2(2\Gamma - \gamma)) y_2 + \gamma^2 z_2^2) y_1 + (\epsilon - 1)y_2 ((2\Gamma - \gamma)\Gamma y_2^2 + \gamma^2 z_2^2).
\]

First, assume that \( p_4 = 0 \). We compute a Gröbner basis of \( I_2 = \langle uy_1 - 1, p_4 \rangle \) for the order \( u \gg z_1 \gg y_1 \gg z_2 \gg y_2 \gg \epsilon > \gamma > \Gamma \), and we find that the last polynomial defining the ideal is \( p_3 \), whose solutions we already studied.

Finally, assume that \( p_4 \neq 0 \) and \( p_5 = 0 \). The discriminant in \( y_2 \) of the coefficient \( a_5(y_2, z_2) \) of \( y_1 \) in \( p_5 \) is
\[
-4a_4(z_2)\gamma^2 z_2^2
\]
where \( a_4(z_2) \) is as above the coefficient of \( z_1 \) in \( p_3 \). The last components of the solutions are given by
\[
\begin{align*}
a_5(y_2, z_2) &\neq 0 \\
y_1 &= \frac{(1-\epsilon)y_2 ((2\Gamma - \gamma)\Gamma y_2^2 + \gamma^2 z_2^2)}{\Gamma (2\epsilon (\epsilon - 2)(\Gamma - \gamma)z_2 + (\epsilon - 1)^2(2\Gamma - \gamma)) y_2 + \gamma^2 z_2^2}
\end{align*}
\]
and
\[
\begin{align*}
a_5(y_2, z_2) &= 0 \\
y_2 &= ((2\Gamma - \gamma)\Gamma y_2^2 + \gamma^2 z_2^2)
\end{align*}
\]
which, as in the case \( y_1 = z_1 = 0 \), is only \( y_2 = z_2 = 0 \) if \( 2\Gamma > \gamma \). This partial solution completes into \( y_1 = y_2 = z_1 = z_2 = 0 \), which was already known. \( \square \)

**Equilibrium positions.**

**Lemma 5.3.** The equilibrium points of \( \dot{X} = DF - D'G \) are all contained in \( \{ D = D' = 0 \} \).

**Proof.** Assume that at some point, either of the determinants \( D \) and \( D' \) is non-zero, this implies that \( F \) and \( G \) are colinear. Since \( F \) and \( G \) form the first two columns of the matrices whose \( D \) and \( D' \) are the respective determinants, \( D = D' = 0 \) at that point. \( \square \)

**Linearization of the system at equilibrium points.** For each of the components of the set of equilibrium points \( \{ D = D' = 0 \} \) found in the previous paragraph, we inspect the behavior of the system in a neighborhood. Namely, for each equilibrium point \( q \), we write
\[
\frac{d}{dt}(q + \delta q) = (DF - D'G)(q) + A(q) \cdot \delta q + R(\delta q).
\]
where \( A = \text{Jac}_q(DF - D'G) \), so that
\[
\frac{d}{dt}(\delta q) = A(q) \cdot \delta q + R(q)(\delta q).
\]
We can compute \( A(q) \) explicitly: Indeed, let \( f = DF - D'G \). Its first derivative is
\[
df(q)(u) = dD(q)(u)F(q) + D(q)dF(q)(u) - dD'(q)(u)G(q) - D'(q)dG(q)(u),
\]
(5.10)
so
\[ A(q) = \nabla D(q) \cdot F(q) + D(q) \text{Jac}_q(F)(q) - \nabla D'(q) \cdot G(q) - D'(q) \text{Jac}_q(G)(q). \]

We examine the eigenvalue decomposition of \( A(q) \).

**Solution 1** (5.5). If \( z_1 = z_2 = \frac{\gamma - 2\Gamma}{\Gamma - \gamma} \), the characteristic polynomial of \( A(q) \) factors as
\[
T^2 \left( T - \gamma^2 (2\Gamma - \gamma)^2 (\epsilon - 1) P_{y_1}^2 \right)^2
\]
The matrix \( A(q) \) is diagonalizable.

**Solution 2** (5.6). If \( y_1 = y_2 = 0 \) and \( z_1 = z_2 \), the characteristic polynomial of \( A(q) \) is
\[
T^4
\]
The Jacobian matrix \( A(q) \) can be trigonalized as
\[ A(q) = P^{-1} \begin{bmatrix} 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{bmatrix} P \]
with the transition matrix
\[
P = \begin{bmatrix}
0 & 0 & 1 & -1 \\
0 & 1 & 0 & 1 \\
\epsilon \gamma^3 (\epsilon - 1) (\epsilon - 2) z_1^2 P_{z_1} & 1 & 0 & 0 \\
\epsilon \gamma^3 (\epsilon - 1) (\epsilon - 2) z_1^2 P_{z_1} & 0 & 0 & 0
\end{bmatrix}.
\]

**Solution 3** (5.7). If \( z_1 = z_2 = \Gamma P_{y_1}^2 (\gamma - 2\Gamma)/2(\Gamma - \gamma)a_3 \), the characteristic polynomial of \( A(q) \) factors as
\[
T^2 \left( T + \frac{b_3}{a_3} \right) \left( T - \frac{b_3}{a_3} \right)
\]
with
\[ b_3 = \Gamma \gamma^2 (\epsilon - 1) P_{y_1} P_{y_2}(2\Gamma - \gamma)^2. \]
The matrix \( A(q) \) is diagonalizable.

**Solution 4** (5.8). If \( y_2 = \frac{y_2 z_2}{(1-\epsilon) z_1} \) and \( z_1 = \frac{(2\Gamma - \gamma) z_2}{a_4} \), the characteristic polynomial of \( A(q) \) factors as
\[
T^2 \left( T - \frac{b_4}{a_4} \right) \left( T + \frac{b_4}{a_4} \right)
\]
with
\[ b_4 = 2\epsilon^2 \gamma^3 z_2^3 (\epsilon - 1)(\epsilon - 2)^2 (2\Gamma - \gamma)(\Gamma - \gamma) P_{z_2}. \]
The matrix \( A(q) \) is diagonalizable.

**Solution 5** (5.9). If \( z_1 = \frac{z_2 y_1}{(1-\epsilon) y_2} \) and \( y_1 = \frac{(1-\epsilon) y_2 ((2\Gamma - \gamma) \Gamma y_2^2 + \gamma^2 z_2^2)}{a_5} \), the characteristic polynomial of \( A(q) \) factors as
\[
T^2 \left( T - \frac{b_4 (\Gamma y_2^2 + \gamma (z_2^2 + z_2)) \Gamma y_2^2}{a_5} \right)^2.
\]
The matrix \( A(q) \) is diagonalizable.
Special points. There are two points at which $A$ vanishes: the North pole $N = (0,0,0,0)$ and $S = (0,z_S,0,z_S)$. Both points are such that $D = D' = 0$, $\nabla D = \nabla D' = 0$, and additionally, at the North pole, $F(N) = 0$.

The North pole is on solutions 2, 3, 4 and 5.

The remainder at $N$ is cubic:

\[
\frac{d}{dt}(N + \delta q) = R(N)(\delta q) = O(\|\delta q\|^3).
\]

The point $S$ is the intersection of solutions 1 and 2. The remainder at $S$ is quadratic.

Higher order studies for the special points.

Quadratic approximation at $S$. We now study the quadratic component $H_2 = Q(S)$ of the remainder $R(S)$:

\[
\frac{d}{dt}(q + \delta q) = (DF - D'G)(q) + A(q)(\delta q) + Q(q)(\delta q) + O(\|\delta q\|^3),
\]

with $\frac{d}{dt}(S) = (DF - D'G)(S) = 0$ and $A(S) = 0$.

We can compute $Q$ by differentiating $f = DF - D'G$ again, as was done in [6, Sec. 3.4]. Differentiating (5.10) along $q$ again, the second derivative of $f$ is

\[
d^2 f(q)(u,v) = d^2 D(q)(u,v)F(q) + dD(q)(u)dF(q)(v) + dD(q)(v)dF(q)(u)
\]
\[\quad - d^2 D'(q)(u,v)G(q) - dD'(q)(u)dG(q)(v) - dD'(q)(v)dG(q)(u)
\]

(5.11)

Note that second derivatives of $F$ and $G$ are 0, since their coordinates are affine in $q$.

We wish to compute $H_2(\delta q) = Q(S)(\delta q, \delta q) = \frac{1}{2}d^2 f(q)(\delta q, \delta q)$. Since $dD(S) = dD'(S) = 0$, we find in the end that

\[
H_2(\delta q) = h_2(\delta q)F(S) - h'_2(\delta q)G(S),
\]

with

\[
F(S) = \begin{pmatrix} 0, \gamma(2\Gamma - \gamma), 0, \gamma(2\Gamma - \gamma) \end{pmatrix}^t,
\]

\[
G(S) = \begin{pmatrix} \frac{1}{2}(2\Gamma - \gamma), 0, (1 - \epsilon)\gamma \end{pmatrix}^t,
\]

\[
h_2(\delta q) = \frac{1}{2}d^2 D(S)(\delta q, \delta q) = (1 - \epsilon)(\delta z_1 - \delta z_2)(\delta z_1 - (1 - \epsilon)^2\delta z_2)(2\Gamma - \gamma)^2,
\]

\[
h'_2(\delta q) = \frac{1}{2}d^2 D'(S)(\delta q, \delta q) = (1 - \epsilon)(\delta z_1 - \delta z_2)(\delta y_2(\epsilon - 1) + \delta y_1)(2\Gamma - \gamma)^2\gamma^2.
\]

Following [6] and [27], we study the projection of the differential equation $\dot{v} = H_2(v)$ on the sphere $S^3$. Let $w = v/\|v\|$ be this projection, it satisfies the differential equation

\[
\dot{w} = \frac{1}{\|v\|^2} \left( \dot{v}\|v\| - v\frac{\langle v, \dot{v} \rangle}{\|v\|} \right)
\]
\[= \frac{H_2(v)}{\|v\|} - \frac{\langle v, H_2(v) \rangle}{\|v\|^3}v
\]
\[= \|v\|(H_2(w) - \langle w, H_2(w) \rangle w)
\]
so we are to study the following differential equation on the sphere $S^3$:

$$\dot{v} = H_2(v) - \langle v, H_2(v) \rangle v =: H_2^2(v).$$

Invariants are related to the eigenvalues of the linearization of $H_2^2$ at points where $H_2^2(v) = 0$. Those points are:

- lines of non-isolated singular points of $H_2$, that is vectors $v$ such that $H_2(v) = 0$
- ray solutions, that is vectors $\xi$ such that there exists $\lambda \in \mathbb{R}\setminus\{0\}$, $H_2(\xi) = \lambda \xi$.

We study the linearization of $H_2^2$ in some neighborhood of these solutions in $S^3$.

**Proposition 5.4.** The blow-up at point $S$ has no ray solution, and two sets of non-isolated singularities:

1. the projective plane $\delta z_1 = \delta z_2$;
2. the projective line $\delta y_2 = (1 - \epsilon)\delta y_1$, $\delta z_1 = (1 - \epsilon)^2 \delta z_2$.

In the first case, the Jacobian of the system is nilpotent. In the second case, it is diagonalizable with non-zero eigenvalues:

$$\frac{1}{2} \left( \delta y_2 + 1 \pm \sqrt{\delta y_2^2 + (2\epsilon - 1)^2(\delta y_2 + 1) - 4(\epsilon - 1)^4 - 2\delta y_2 + 1} \right).$$

**Proof.** First we study ray solutions. Let $\xi$ be a vector on a ray solution, such that $H_2(\xi) = \lambda \xi$.

Let $\alpha \xi$ be another vector on the same line ($\alpha \in \mathbb{R}$), since $H_2$ is homogeneous with degree 2, one has

$$H_2(\alpha \xi) = \alpha^2 H_2(\xi) = \alpha^2 \lambda \xi = \alpha \lambda (\alpha \xi).$$

So each line or ray solutions contains a unique $\xi_0$ such that $H_2(\xi_0) = \xi_0$.

A Gröbner basis of the system $\langle H_2(\delta q) - \delta q \rangle$ is given by $\{\delta y_1, \delta z_1, \delta y_2, \delta z_2\}$, so there is no non-trivial ray.

This can also be seen in the following way: let $\delta q$ be a vector such that $H_2(\delta q) = \delta q$. By the structure of the vector $F(S)$, $\delta q$ satisfies $\delta z_1 = \delta z_2$, and so $h_2(\delta q) = h'_2(\delta q) = 0$, so $H_2(\delta q) = 0$, which, by hypothesis, implies that $\delta q = 0$.

We now consider non-isolated singular points of $H_2$, that is the zeroes of $H_2$. Since $F(S)$ and $G(S)$ are linearly independent, those points are exactly the zeroes of $h_2$ and $h'_2$, as described in the statement of the proposition.

Then we study the linearization of $H_2^2$ in some neighborhood of these solutions in $S^3$. First we consider vectors $\delta q$ such that $\delta z_1 = \delta z_2$, we may perform the computations in the affine chart given by $\delta z_1 \neq 0$, with coordinates $\delta y_1 = \delta y_1/\delta z_1$, $\delta y_2 = \delta y_2/\delta z_1$, $\delta z_2 = \delta z_2/\delta z_1$. The differential equation becomes

$$\frac{d}{dt} \begin{pmatrix} \delta y_1 \\ \delta y_2 \\ \delta z_1 \end{pmatrix} = \delta z_2 C(\delta y_1, \delta y_2, \delta z_1)$$

with $C$ a polynomial vector field of degree 3.

At $\delta z_1 = 1$, its Jacobian is nilpotent:

$$\begin{pmatrix} 0 & 0 & (\epsilon - 1)(\delta y_1(\epsilon(\epsilon - 1) - 1) + \delta y_2(1 - \epsilon))(2\Gamma - \gamma)^2 \gamma^3 \\ 0 & 0 & (\epsilon - 1)(\delta y_1(2\epsilon(\epsilon - 2) + 1) + \delta y_1(\epsilon - 1))(2\Gamma - \gamma)^2 \gamma^3 \\ 0 & 0 & 0 \end{pmatrix}$$
Then we consider vectors $\delta q$ such that $\delta y_2 = (1 - \epsilon)\delta y_1$, $\delta z_1 = (1 - \epsilon)^2 \delta z_2$. This time we use the chart $\delta z_2 \neq 0$ with coordinates $\delta y_1 = \delta y_1 / \delta z_2$, $\delta y_2 = \delta y_2 / \delta z_2$, $\delta z_1 = \delta z_1 / \delta z_2$. As above, we compute the differential equation in this chart, linearize the resulting vector field, and evaluate this Jacobian at $\delta y_2 = (1 - \epsilon)\delta y_1$ and $\delta z_1 = (1 - \epsilon)^2$. This matrix has rank 2 and is diagonalizable with non-zero eigenvalues:

$$
\frac{1}{2} \left( \delta y_2 + 1 \pm \sqrt{\delta y_2^2 + (2\epsilon - 1)^2(\delta y_2 + 1) - 4(\epsilon - 1)^4 - 2\delta y_2 + 1} \right).
$$

\[\square\]

**Cubic approximation at $N$.** We perform the same study at the North pole $N$. With expression (5.11), we can verify that the quadratic component of $R(N)$ is 0. Indeed, $F(N) = 0$ and $d^2D'(N) = 0$.

Further differentiating along $q$, we obtain

$$
H_3(\delta q) := \frac{1}{6} d^3 f(N)(\delta q, \delta q, \delta q) = \frac{1}{6} \left( 3d^2 D(N)(\delta q, \delta q) F(\delta q) - d^3 D'(N)(\delta q, \delta q, \delta q) G(N) \right)
$$

Note that since we centered the coordinates at the North pole, $F$ is linear in $q$, so $dF(q) = F$, and $G$ is affine in $q$, so $dG(q)$ is constant.

As in the previous subsection, we study the projection of the differential equation $\dot{v} = H_3(v)$ on the sphere $S^3$, and its equilibrium points, which form lines of non-isolated singular points and ray solutions.

**Proposition 5.5.** The cubic blow-up at the North pole $N$, for admissible values of the parameters, has two sets of ray solutions:

1. the projective line

$$
\begin{cases}
\delta z_1 = \delta z_2 = 0 \\
(\epsilon - 1)\delta y_1 + \delta y_2 = \frac{1}{\Gamma(2\Gamma - \gamma)\sqrt{1 - \epsilon}};
\end{cases}
$$

(5.12)

2. the quadric

$$
\begin{cases}
\delta y_1 = \delta y_2 = 0 \\
((\epsilon - 1)\delta z_1 - \delta z_2)^2 - (\epsilon - 2)\delta z_1 \delta z_2 = \frac{1}{\gamma^3(2\Gamma - \gamma)(1 - \epsilon)}.
\end{cases}
$$

and three sets of real non-isolated singularities:

1. the plane

$$
\begin{cases}
\delta z_1 = \delta z_2 \\
\delta y_1 (1 - \epsilon) = \delta y_2
\end{cases}
$$

(5.13)

2. the plane

$$
\begin{cases}
\delta y_2 = (1 - \epsilon)\delta y_1 \\
\delta z_2 = (1 - \epsilon)^2 \delta z_1
\end{cases}
$$

(5.14)
(3) the surface defined by
\[
\begin{align*}
0 &= \Gamma(\epsilon - 1)(2\Gamma - \gamma)\delta y_1^2 + (2\Gamma - \gamma)\Gamma\delta y_1 \delta y_2 + \gamma^2(\epsilon - 1)\delta z_1^2 - \gamma^2(\epsilon - 1)\delta z_2^2 \\
0 &= \Gamma(\epsilon - 1)(2\Gamma - \gamma)\delta y_1 \delta y_2 - \gamma^2\delta z_1\delta z_2 + (2\Gamma - \gamma)\Gamma\delta y_2^2 + \gamma^2\delta z_2^2 \\
0 &= \delta y_1 \delta z_2 + (\epsilon - 1)\delta y_2 \delta z_1.
\end{align*}
\]
(5.15)

For points on the line (5.12), the linearization of $H_3^2$ is diagonal: the vectors $(1, 0, 0)$ and $(0, 0, 1)$ are eigenvectors, with the same eigenvalue, and the vector $(0, 0, 1)$ is in the kernel.

For isolated singularities satisfying (5.13), the matrix is not diagonalizable, its Jordan form has the following structure:
\[
\begin{bmatrix}
0 & 0 & 0 \\
0 & * & 1 \\
0 & 0 & *
\end{bmatrix}
\]

For isolated singularities satisfying (5.14), the matrix is diagonalizable with 3 non-zero eigenvalues.

Proof. First we study ray solutions. Let $\xi$ be a vector on a ray solution, such that $H_3(\xi) = \lambda \xi$.

Let $\alpha \xi$ be another vector on the same line ($\alpha \in \mathbb{R}$), one has $H_3(\alpha \xi) = \alpha^3 H_3(\xi) = \alpha^3 \lambda \xi = \alpha^2 \lambda (\alpha \xi)$.

So unlike in the quadratic case, a line of ray solutions contains 2 vectors $\xi_1, \xi_2$ such that either $H_3(\xi_1) = H_3(\xi_2) = 1$ or $H_3(\xi_1) = H_3(\xi_2) = -1$.

In order to study ray solutions, we compute a Gröbner basis of the system $H_3(\delta q) - \iota \delta q = 0, \iota^2 = 1$, for an order eliminating $\alpha$. We find that the basis contains
\[
\{\delta y_1 \delta z_1, \delta y_1 \delta z_2, \delta y_2 \delta z_1, \delta y_2 \delta z_2\}
\]
so either $\delta y_1 = \delta y_2 = 0$ or $\delta z_1 = \delta z_2 = 0$.

If $\delta z_1 = \delta z_2 = 0$, computing a new Gröbner basis of the system, saturating with $\Gamma - \gamma$ and $\delta y_1$ shows that $\delta y_1$ and $\delta y_2$ must satisfy
\[
\left((\Gamma^2(\epsilon - 1)(2\Gamma - \gamma)^2(\epsilon - 1)\delta y_1 + \delta y_2)^2\right)^2 - 1 = 0.
\]

Since $\epsilon - 1 < 0$, $2\Gamma - \gamma > 0$ and $\Gamma > 0$, this defines 2 lines of real solutions given by
\[
(\epsilon - 1)\delta y_1 + \delta y_2 = \pm \frac{1}{\Gamma(2\Gamma - \gamma)\sqrt{\Gamma - \epsilon}}.
\]
Those lines are equivalent in the projective space: each line of ray-solutions contains a vector in both lines.

If $\delta y_1 = \delta y_2 = 0$, the same technique shows that $\delta z_1$ and $\delta z_2$ must satisfy
\[
\left((\gamma^3(2\Gamma - \gamma)(\epsilon - 1)
\left(((\epsilon - 1)\delta z_1 - \delta z_2)^2 - (\epsilon - 2)\delta z_1 \delta z_2\right)\right)^2 - 1 = 0,
\]
which defines 1 quadric
\[ ((\epsilon - 1)\delta z_1 - \delta z_2)^2 - (\epsilon - 2)\delta z_1 \delta z_2 = \frac{1}{\gamma^3 (2\Gamma - \gamma)(1 - \epsilon)}. \]

We now consider non-isolated singularities, that is zeros of \( H_3 \). To this end, we compute a Gröbner basis of the system \( H_3(\delta y) = 0 \), saturating by \( \gamma - \Gamma, \gamma, \Gamma, \epsilon - 1 \) and \( 2\Gamma - \gamma \). Factoring the results, it appears that the solutions split into 5 cases:

1. \( \delta z_1 = \delta z_2 \)
2. \( \delta z_1 = -\delta z_2 \)
3. \( \delta z_1 = 0 \)
4. \( \delta z_2 = (1 - \epsilon)^2 \delta z_1 \)
5. otherwise.

In the case 1, a Gröbner basis is given by
\[
\begin{align*}
& \{ \delta y_1(\delta y_1(\epsilon - 1) + \delta y_2)^2 \\
& \delta y_2(\delta y_1(\epsilon - 1) + \delta y_2)^2 \\
& \delta z_2(\delta y_1(\epsilon - 1) + \delta y_2)^2 \\
& \delta z_1 - \delta z_2 \}
\end{align*}
\]
and the solutions form the plane (5.13).

In the case 2, saturating by \( \delta z_1 - \delta z_2 \), a Gröbner basis is given by
\[
\left\{ \delta y_1 + (1 - \epsilon)\delta y_2, \delta z_1 + \delta z_2, \Gamma(\epsilon^2 + (2 - 2\epsilon)(2\gamma - \gamma)\delta y_2^2 + 2\gamma^2 \delta z_2^2) \right\}
\]
which has no real non-zero solution for admissible values of the parameters.

In the case 3, saturating by \( \delta z_1 - \delta z_2 \) and \( \delta z_1 + \delta z_2 \), a Gröbner basis is given by
\[
\{ \delta y_1, \delta z_1, (2\Gamma - \gamma)\Gamma\delta y_2^2 + \gamma^2 \delta z_2^2 \}
\]
which has no real non-zero solution for admissible values of the parameters.

In the case 4, saturating by \( \delta z_1 \pm \delta z_2 \) and \( \delta z_1 \), a Gröbner basis is given by
\[
\{ \delta y_1 \delta z_2 + (\epsilon - 1)\delta y_2 \delta z_1, \delta y_2 + (\epsilon - 1)\delta y_1, \delta z_2 - (\epsilon - 1)^2 \delta z_1 \}
\]
and the solutions form the plane (5.14).

Finally, for the case 5, we compute a Gröbner basis, saturating by all the previous conditions. This basis is
\[
\left\{ \Gamma(\epsilon - 1)(2\Gamma - \gamma)\delta y_1^2 + (2\Gamma - \gamma)\Gamma \delta y_1 \delta y_2 + \gamma^2 (\epsilon - 1)\delta z_1^2 - \gamma^2 (\epsilon - 1)\delta z_1 \delta z_2 \\
\Gamma(\epsilon - 1)(2\Gamma - \gamma)\delta y_1 \delta y_2 - \gamma^2 \delta z_1 \delta z_2 + (2\Gamma - \gamma)\Gamma \delta y_2^2 + \gamma^2 \delta z_2^2 \\
\delta y_1 \delta z_2 + (\epsilon - 1)\delta y_2 \delta z_1 \\
(\Gamma(2\Gamma - \gamma)(\epsilon - 1)^2 \delta y_2^2 + \delta z_2^2 \gamma^2) \delta z_1 - (2\Gamma - \gamma)\Gamma \delta z_2 \delta y_2^2 - \gamma^2 \delta z_2^3 \right\}
\]
The fourth polynomial is a combination of the other 3, and the solutions form the surface (5.15).

For the second part of the proposition, as in the quadratic case, we study the linearization of \( H_3^\sharp : \dot{v} = H_3(v) - \langle v, H_3(v) \rangle v \). In the affine chart with \( \delta y_1 \neq 0 \), with coordinates \( \delta z_1 = \delta z_1/\delta y_1, \delta y_2 = \delta y_2/\delta y_1 \) and \( \delta z_2 = \delta z_2/\delta y_1 \), the differential
equation $\dot{v} = H_3(v)$ becomes

$$\frac{d}{dt} \begin{pmatrix} \delta z_1 \\ \delta y_2 \\ \delta z_2 \end{pmatrix} = \delta y_2^2 \bar{Q}(\delta z_1, \delta y_2, \delta z_2)$$

with $\bar{Q}$ a polynomial vector field of degree 4. In this chart, $H_3^x$ becomes $\dot{v} = \bar{Q}(v)$.

We conclude by evaluating the Jacobian of $\bar{Q}$ at the relevant points. □

5.3.2. General case.

Singularities of $\{D = H_G = \{H_G, H_F\} = 0\}$.

Proposition 5.6. The set of singularities of $\{D = H_G = \{H_G, H_F\} = 0\}$ is given, generically on authorized values of $\gamma, \Gamma$, by

1. the plane $z_1 = z_2 = z_S$;
2. the line $z_1 = z_2, y_1 = y_2 = 0$;
3. an irreducible variety of dimension 5.

If $\gamma = \Gamma$, solution 2 becomes a surface defined by $z_1 = z_2, y_2 = (1 - \epsilon)y_1$.

Proof. By definition of $D = \{\{H_G, H_F\}, H_G\}$, we want to study the zeroes of

$$\begin{cases} 0 = p \cdot G \\ 0 = p \cdot [G, F] \\ 0 = p \cdot [[G, F], G] \end{cases}$$

(D)

The singularities of this variety is the set of points at which the matrix

$$\begin{bmatrix} G & [G, F] & [[G, F], G] \\ v_1 & v_2 & v_3 \end{bmatrix}$$

has rank at most 2.

We encode that with the incidence variety

$$\begin{bmatrix} G & [G, F] & [[G, F], G] \\ v_1 & v_2 & v_3 \end{bmatrix} \cdot \begin{bmatrix} L_1 \\ L_2 \\ L_3 \end{bmatrix} = \begin{bmatrix} 0 \\ 0 \\ 0 \\ 1 \end{bmatrix}$$

with new variables $L = L_1, L_2, L_3$ and random numbers $v_1, v_2, v_3$. This gives us a system of 4 polynomial equations in the 10 unknowns $y_1, z_1, z_2, L_1, L_2, L_3, \Gamma, \gamma, \epsilon$. We eliminate $L_1, L_2, L_3$ from the ideal in order to recover the projection, and we saturate by $1 - \epsilon$.

We then compute a Gröbner basis for the elimination order $y_1 > y_2 > z_1 > z_2 > \gamma > \Gamma > \epsilon$, this basis contains 10 polynomials, some of which have factors with multiplicity greater than 1 or are divisible by $1 - \epsilon$ or $\gamma$. We take the square-free form of this basis, and we saturate by $1 - \epsilon$ and $\gamma$ before computing a new Gröbner basis for the same order. The result is a set of 11 polynomials which includes

$$(z_1 - z_2) ((\epsilon - 1)P_{z_2} y_1 + P_{z_1} y_2).$$
First, we add \( z_1 - z_2 \) to the ideal. Once again, we compute a Gröbner basis for the elimination order \( y_1 > y_2 > z_1 > z_2 \gg \gamma > \Gamma > \epsilon \), take its square-free form, and recompute a Gröbner basis. The result contains the polynomial \( P_{z_2}P_{y_2} \).

The solutions decompose into 2 algebraic sets, defined by

\[
z_1 = z_2 = z_S = \frac{2\Gamma - \gamma}{2\Gamma - 2\gamma}
\]

and (adding \((\epsilon - 1)y_1 + y_2\) to the ideal and saturating by \(P_{z_2}, \epsilon\) and \(\epsilon - 2\))

\[
\begin{aligned}
0 &= y_1(\Gamma - \gamma) \\
0 &= y_2(\Gamma - \gamma) \\
0 &= P_{y_2} = (\epsilon - 1)y_1 + y_2 \\
0 &= z_1 - z_2.
\end{aligned}
\]

The latter is generically (i.e. if \( \Gamma \neq \gamma \)) defined by

\[
\begin{aligned}
y_1 &= y_2 = 0 \\
z_1 &= z_2
\end{aligned}
\]

or, if \( \Gamma = \gamma \) (that is if the spin we consider is water), by

\[
\begin{aligned}
y_1 &= \frac{y_2}{\epsilon - 1} \\
z_1 &= z_2.
\end{aligned}
\]

Then, starting again with the whole ideal, we add \((\epsilon - 1)P_{z_2}y_1 + P_{z_2}y_2\) to the ideal and saturate by \(z_1 - z_2\). The result, generically on \((\Gamma, \gamma, \epsilon)\), is an irreducible surface. □

**Locus of** \( \{D = D' = H_G = \{H_G, H_F\} = 0\} \).

**Proposition 5.7.** The solutions form the union of the hyperplane defined by

\[
z_1 = z_2
\]

and the hypersurface

\[
y_1 = -\frac{y_2P_{z_1}}{(\epsilon - 1)P_{z_2}}.
\]

**Proof.** Points such that \( \{D = D' = H_G = \{H_G, H_F\} = 0\} \) satisfy

\[
\begin{aligned}
0 &= p \cdot G \\
0 &= p \cdot [G, F] \\
0 &= p \cdot [[G, F], G] \\
0 &= p \cdot [[G, F], F]
\end{aligned}
\]

The projection of these points onto the space \((y_1, z_1, y_2, z_2)\) is given by the vanishing of the determinant \(\Delta'\), defined as

\[
\Delta' = \det \begin{bmatrix}
(1 - \epsilon)G & [(1 - \epsilon)G, F] & [(1 - \epsilon)G, (1 - \epsilon)G] & [(1 - \epsilon)(G, F), F] \\
\end{bmatrix}
\]

\[
= (1 - \epsilon)^2 \det \begin{bmatrix}
\end{bmatrix}
\]
This determinant factors as
\[ -2 (\epsilon - 1)^2 (2 \gamma - 1)(\Gamma - 1)(z_1 - z_2) \left[ (\epsilon - 1) P_{z_1} y_1 + P_{z_2} y_2 \right] \]

Equilibrium positions.

Lemma 5.8. All equilibrium points of \( \dot{Z} = D\vec{H}_F - D'\vec{H}_G \) satisfying \( H_G = \{ H_G, H_F \} = 0 \) are contained in \( \{ D = D' = 0 \} \).

Proof. Recall that \( \vec{H}_F \) is defined as
\[ \vec{H}_F = \begin{bmatrix} \frac{\partial H_F}{\partial p} \\ -\frac{\partial F}{\partial q} \end{bmatrix} = \begin{bmatrix} F - p \cdot \frac{\partial H_F}{\partial q} \\ -p \cdot \frac{\partial H_F}{\partial q} \end{bmatrix} \]
and \( \vec{H}_G \) is defined in the same way. Let \( z = (q, p) \) be a point such that
\[ D(z)\vec{H}_F(z) - D'(z)\vec{H}_G(z) = 0, \]
by looking at the first 4 components of this system, we see that the vectors \( F(q) \) and \( G(q) \) are colinear.

Introduce new variables \( X_D \) and \( X_D' \) and consider the ideal generated by
- \( X_D F(q) - X_D' G(q) = 0 \)
- \( H_G(z) = \{ H_G, H_F \} (z) = 0 \)
- \( D(z)\vec{H}_F(z) - D'(z)\vec{H}_G(z) = 0 \)
saturated by \( X_D \) and \( X_D' \). Computing a Gröbner basis of this ideal (for any order) yields that this ideal is actually \( \langle 1 \rangle \), and so the associated system has no solution.

Hence, at an equilibrium point, either \( D(z) \) or \( D'(z) \) has to be 0. Since both \( O \) (the center of the Bloch ball) and \( N \) (the north pole) are on the hyperplane \( z_1 = z_2 = 0 \), which is contained in \( \{ D = D' = 0 \} \), we may assume that the point \( z \) is neither \( O \) nor \( N \). Hence, \( F(z) \) and \( G(z) \) are non-zero, and so \( D(z) = D'(z) = 0 \).

Eigenvalues of the linearization. We consider the eigenvalue decomposition of the matrix
\[ \mathcal{A} = \text{Jac}(D\vec{H}_F - D'\vec{H}_G) \]
on equilibrium point, given as the union of points satisfying Eq. (5.16) and (5.17).

Solutions of Eq. (5.16). If \( z_1 = z_2 \), the matrix \( \mathcal{A} \) has rank 2, and its characteristic polynomial is
\[ T^6 \left( T^2 - \frac{\epsilon(\epsilon - 1)(\epsilon - 2)}{2P_{y_1} (\Gamma - \gamma)} \left( 8(\Gamma - \gamma)^2 P_{y_1} y_1 y_2 + P_{z_2} (2P_{z_2} - \gamma) P_{y_2} \right) \right) \]
\[ + \left( \frac{(\epsilon - 1)\gamma P_{y_1} (2\Gamma - \gamma)}{y_1} \right)^2 \]

Solutions of Eq. (5.17). If \( y_2 = -\frac{y_2 P_{y_1}}{(\epsilon - 1) P_{z_2}} \), the matrix \( \mathcal{A} \) has rank 2, and its characteristic polynomial is
\[ T^6 \left( T^2 - 4\epsilon y_2 (\epsilon - 1)(\epsilon - 2)(\Gamma - \gamma) T + \left( \frac{2\gamma (z_1 - z_2)(\epsilon - 1)^2 (2\Gamma - \gamma)(\Gamma - \gamma)}{P_{z_1}} \right)^2 \right). \]
The discriminant of the degree 2 factor factors as
\[ 16(\Gamma - \gamma)^2(\epsilon - 1)^2 (a_6(z_1, y_2) - b_6(z_1, z_2)) (a_6(z_1, y_2) + b_6(z_1, z_2)) \]
with
\[ a_6(z_1, y_2) = \epsilon(\epsilon - 2)P_{z_1}y_2 \]
\[ b_6(z_1, z_2) = (2\Gamma - \gamma)(\epsilon - 1)\gamma(z_1 - z_2) \]
which induces the following classification of the eigenvalues of \( A \):
- if \(|a(z_1, y_2)| > |b(z_1, z_2)|\): 2 single real eigenvalues;
- if \(|a(z_1, y_2)| = |b(z_1, z_2)|\): 1 double real eigenvalue;
- if \(|a(z_1, y_2)| < |b(z_1, z_2)|\): 2 single complex eigenvalues.

6. Conclusion

In this article we made a complete use of the state of the art in geometric, symbolic and numeric techniques to analyze the problem of saturating a pair of spins in relation with the \( B_1 \)-inhomogeneities, that is inhomogeneities of the applied RF-field. We extend the results in many directions. First of all, the time minimal syntheses for a single spin are classified, taking into account the relaxation parameters and the control bounds. For a pair of spins, the crucial theoretical problem is to classify the singularities of the extremal flow. This is realized using symbolic computations based on Gröbner basis to compute the singularities and linear or quadratic, cubic approximations of some crucial frame points. This is not sufficient to make a topological classification of the behaviours, since Grobman-Hartman cannot be applied in general for non isolated singularities, see [6]. Nevertheless numerical methods using continuation techniques can be used to analyze the singularities. One feature of the problem is the existence of many local optima and applications of LMI techniques are particularly important to compare the different local optima obtained using direct or indirect numerical schemes implemented in the Bocop and HamPath software. Hence, we believe that this article complete in many directions the results and techniques obtained in previous articles. It is a relevant step in the problem of determining the cartography of the global optima with respect to the relaxation parameters in the (ideal) contrast problem in MRI and to provide substantial improvements in existing software in MRI.

Also in the contrast of geometric optimal control it is a significant step to handle complex 4-D problems.

References


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