# Algebraic-geometric techniques for the feedback classification and robustness of the optimal control of a pair of Bloch equations with application to magnetic resonance imaging

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The aim of this article is to classify using computer algebraic methods the singular trajectories of a pair of controlled Bloch equations in view of analysing the robustness with respect to the parameters and  $B_1$ -inhomogeneity of the optimal solution to the contrast and multisaturation problems in magnetic resonance imaging. They are completed by preliminary numerical simulations.<sup>1</sup>

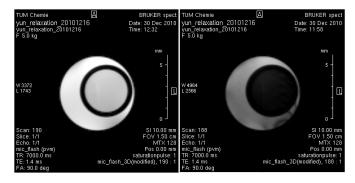
# 1 Introduction

Optimal control algorithms were introduced in Nuclear Magnetic Resonance (NMR) to improve the control field very recently [13] and at the end of the nineties, new methods appeared in optimal control of NMR systems both from the analytical and numerical points of view [19, 34]. More recently, under the impulse of S. Glaser, combination of geometrical optimal control based on the Maximum Principle [30] and related numerical algorithms: gradient methods [20], shooting and continuation methods [14] lead to sophisticated results starting from a complete solution to the time-minimal saturation of a single spin [23], and applications to the contrast problem in magnetic resonance imagery (MRI), see [5, 22].

They are the basis to the numeric computations of robust optimal controls with respect to  $B_0$  and  $B_1$  inhomogeneities and were validated vert recently by *in vitro* (Figure 1) and *in vivo* experiments [8, 33, 35] (Figure 2).

The mathematical model which very accurately is suitable for analyzing such problems is to consider an ensemble of spins, each spin being described by a magnetization vector M =

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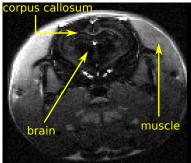


Figure 1: Experimental results: The inner circle shape Figure 2: Contrast optimization sample mimics the deoxygenated blood, where  $T_1 = 1.3$ s and  $T_2 = 50$ ms; the outside moon shape sample corresponds to the oxygenated blood, where  $T_1 = 1.3$ s and  $T_2 = 200$ ms. (left) Without control, (right) Optimized contrast.

in a in vivo setting. Species: brain - parietal muscle

 $(M_X, M_Y, M_Z)$  in the laboratory frame 0XYZ whose evolution satisfies the so-called Bloch equation

$$\frac{\mathrm{d}M}{\mathrm{d}t} = \gamma M \wedge B + R(M),\tag{1}$$

where  $\gamma$  is the gyromagnetic ratio, B(t) is the total magnetic field applied to the system which decomposes into

$$B(t) = B_0 + B_1(t)$$

where  $B_0$  is a strong constant magnetic field oriented along the Z axis, and  $B_1(t)$  is the control RF-field in the transverse (X,Y) plane. The R(M) term represents the dissipation, of the form

$$R(M) = \left(\frac{M_X}{T_2}, \frac{M_Y}{T_2}, \frac{M_Z - M_0}{T_1}\right)$$

where  $M_0$  is the equilibrium magnetization, which can be normalized to 1 using the rescaling  $M \mapsto M/M_0$ , and  $T_1, T_2$  are the relaxation parameters which are the chemical signatures of the observed species. The control components are denoted  $u(t) = -\gamma B_Y$ ,  $v(t) = -\gamma B_X$  and up to a time rescaling, one can impose the control bounds:  $u^2 + v^2 = 1$ .

The Bloch equations can be written in a rotating frame 0xyz with M(t) = S(t)q(t), q = (x, y, z),  $S(t) = \exp(\omega t \Omega_z), \, \Omega_z = \begin{bmatrix} 0 & -1 & 0 \\ 1 & 0 & 0 \\ 0 & 0 & 0 \end{bmatrix}$  as:

$$\frac{\mathrm{d}}{\mathrm{d}t} \begin{bmatrix} x \\ y \\ z \end{bmatrix} = \begin{bmatrix} -1/T_2 & -\Delta \omega & u_2 \\ \Delta \omega & -1/T_2 & -u_1 \\ u_2 & u_1 & -1/T_1 \end{bmatrix} \begin{bmatrix} x \\ y \\ z \end{bmatrix} + \begin{bmatrix} 0 \\ 0 \\ 1/T_1 \end{bmatrix}$$
 (2)

where  $\Delta\omega=\omega_0-\omega$  is the resonance offset,  $\omega_0=-\gamma B_0$  is the resonance frequency and the

RF-control field is represented as

$$\begin{cases} u_2 = u\cos(\omega t) - v\sin(\omega t) \\ u_1 = u\sin(\omega t) + v\cos(\omega t) \end{cases}$$
(3)

which preserves the control bound  $u_1^2 + u_2^2 \le 1$ .

Finally, in the moving frame the system takes the normalized form

$$\begin{cases} \frac{\mathrm{d}x}{\mathrm{d}t} = -\Gamma x - \Delta \omega y + u_2 z \\ \frac{\mathrm{d}y}{\mathrm{d}t} = \Delta \omega x - \Gamma y - u_1 z \\ \frac{\mathrm{d}z}{\mathrm{d}t} = \gamma (1 - z) + u_1 y - u_2 z. \end{cases}$$
(4)

The resonant case is when  $\Delta \omega = 0$  which will be assumed in the sequel of this article. The system is then written as

$$\frac{\mathrm{d}q}{\mathrm{d}t} = F(q) + u_1 G_1(q) + u_2 G_2(q), u_1^2 + u_2^2 \le 1$$

Thanks to the symmetry of revolution along the Z axis related to the  $B_0$ -polarization, in many cases it is physically relevant to consider the situation with  $u_2 = 0$  and the spin system is a 2-dimensional system, with q = (y, z) and it is simply written as

$$\frac{dq}{dt} = F(q) + uG(q), q = (y, z), |u| \le 1.$$
 (5)

Also note that the Bloch ball  $|q| \le 1$  is invariant for the dynamics provided that the parameters satisfy the physical constraint  $2\Gamma \ge \gamma \ge 0$ .

Our study will concern the following optimal control problems.

First of all, consider a single spin system described by (5). The *saturation problem* is to drive the magnetization vector from the north pole of the Bloch ball N = (0,1) to the center O = (0,0). A major result in NMR was to compute the time-minimal solution, using geometric optimal control based on the maximum principle [21].

According to this principle, a time-minimal solution is given by the equations:

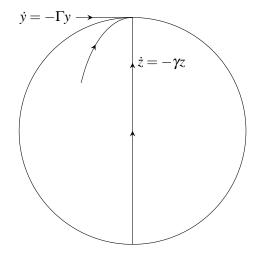
$$\dot{z} = \vec{H}_F(z) + u\vec{H}_G(z), z = (q, p)$$
 (6)

 $H_F$  and  $H_G$  being the Hamiltonians  $p \cdot F(q)$  and  $p \cdot G(q)$ , and moreover the optimal control is given by the maximization condition

$$H_F(z) + uH_G(z) = \max_{|v| < 1} H_F(z) + vH_G(z)$$
(7)

and it is a concatenation of bang arcs  $\delta_+$ ,  $\delta_-$  where u=+1 or u=-1 and the so-called singular arcs where they are obtained by solving  $H_G(z(t))=0$ .

A straightforward computation shows that singular arcs are located on the set where G and [G,F] are colinear and are formed by the two lines y=0 (axis of revolution) and  $z_0=-\frac{\gamma}{2(\Gamma-\gamma)}$  and the interesting situation is when  $2\Gamma>3\gamma$  so that this horizontal line intersects the Bloch ball. More precisely, following [21] we recall the following:



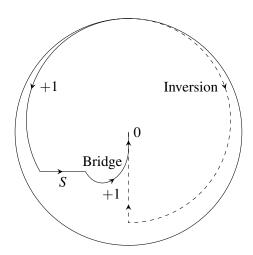


Figure 3: Dynamics under the control  $u \equiv 0$ 

Figure 4: Time-minimal saturation: left:  $2\Gamma > 3\gamma$ , right:  $2\Gamma \le 3\gamma$ 

**Proposition 1.** Provided that  $2\Gamma > 3\gamma$ , the time-minimal solution to the saturation problem in the experimental situation is of the form  $\delta_+\delta_{SH}\delta_+^B\delta_{SV}$  where  $\delta_{SV}$  and  $\delta_{SH}$  are vertical and horizontal singular arcs and the intermediate arc  $\delta_+^B$  is a bang arc connecting the two lines, called a bridge and is related to the property that the singular control as calculated along the singular line is such that  $|u_S| \to \infty$  when  $y \to 0$ .

(Note: in experimental situation, the line  $(\delta_{SH})$  is accessible from the North pole).

All the information about this study is reported on Fig. 4. Note that in the case  $2\Gamma \le 3\gamma$ , the standard inversion sequence  $\delta_+\delta_{SV}$  is time-minimal.

In MRI, we attribute to |q| a level of gray such that |q|=1 corresponds to while and saturation |q|=0 to black. This led to consider the following optimal problems, taking a pair of spins, each governed by (5), with respective parameters  $(\Gamma_1, \gamma_2)$ ,  $(\Gamma_2, \gamma_2)$  and controlled by the same RF-field u (coupling the dynamics).

Problem 1. Saturation of both spins in minimum time. Note that another problem is to saturate a couple of spin systems with the same parameters but taking into account  $B_1$ -inhomogeneity and this leads to consider a pair of systems (5):

$$\begin{cases} \frac{\mathrm{d}q_1}{\mathrm{d}t} = F(q_1) + uG(q_1) \\ \frac{\mathrm{d}q_2}{\mathrm{d}t} = F(q_2) + u(1 - \varepsilon)G(q_2) \end{cases} |u| \le 1$$

where  $\varepsilon$  is a small rescaling parameter related to the control field inhomogeneity.

Problem 2. The second problem which plays a central role in MRI is the contrast problem. One must separate in a fixed transfer time  $t_f$  the contrast between the two species. This amounts to introduce a cost function

$$c(q(t_f)) = \alpha |q_1(t_f)|^2 - \beta |q_2(t_f)|^2$$

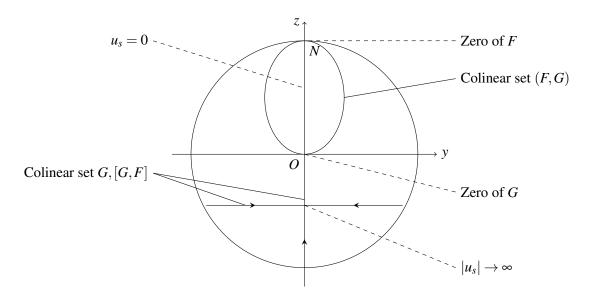


Figure 5: Algebraic sets involved in saturation singular arcs

where  $\alpha$ ,  $\beta$  are weight parameters. The *contrast by saturation* is to saturate one spin, *e.g.* spin 1, imposing  $q_1(t_f) = 0$  while maximizing  $|q_2(t_f)|^2$ .

Both problems correspond to *Mayer problems in optimal control*, that is an optimal problem of the form  $\min_{u(\bullet)} c(q(t_f))$  subject to  $\dot{q} = F + uG$ ,  $|u| \le 1$ , with some prescribed boundary conditions  $q(t_f) \in M$ , with M the terminal manifold defined by f(q) = 0. Again the maximum principle described by (6), (7) applies and is complemented by *transversality conditions* depending upon the cost function and boundary conditions and written as:

$$f(q(t_f)) = 0, p(t_f) = p_0 \frac{\partial c}{\partial a}(q(t_f)) + \delta \frac{\partial f}{\partial a}(q(t_f)), p_0 \le 0, \delta: \text{ constant vector}$$
 (8)

Clearly, the optimal control problems boil down to analyze the so-called singular extremals given by

$$\dot{z} = H_F(z) + uH_G(z), H_G = 0$$
 (9)

where u has to be admissible, that is  $|u| \le 1$ . o The surface  $\Sigma : H_G(z) = 0$  is called the *switching surface*. Computation of the singular extremals can be made explicit. If F, G are two vector fields, the Lie bracket is defined with the convention:

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

and the Poisson bracket of  $H_F$ ,  $H_G$  is given by

$$\{H_F, H_G\}(z) = dH_F(\vec{H}_G(z)) = p \cdot [F, G](q).$$

Hence differentiating  $t \mapsto H_G(z(t)) = 0$ , one gets

$$z \in \Sigma' : H_F(z) = \{H_F, H_G\}(z) = 0$$
 (10)

and the singular control is defined solving

$$\{\{H_G, H_F\}, H_F\}(z) + u\{\{H_G, H_F\}, H_G\} = 0. \tag{11}$$

If F, G are real analytic, this defines a meromorphic equation of the form

$$\frac{\mathrm{d}z}{\mathrm{d}t} = \frac{X(z)}{g(z)}, z \in \Sigma'. \tag{12}$$

Since F and G are affine vector fields and depend linearly on the parameters, this leads to algebraic computations which are presented in this article and performed using computer algebra algorithms. More precisely, they are related to the following problem.

Find rational invariants with respect to the parameters  $((\Gamma_1, \gamma_1), (\Gamma_2, \gamma_2))$  related to the classification of solutions of (12).

Clearly this algebraic study is the first step which complemented by numerical simulations using shooting technique, implemented in the HamPath software and LMI techniques to evaluate the global optimum implemented in the Globoptipoly [18] to classify the optimal strategies and analyse their robustness with respect to the physical parameters [6]. The organization of this article is the following. In Section 2, we present based on [2] the geometric frame relating the classification of singular extremals with the optimal control problems and the action of the feedback group on the system. They are used to generate rational invariants. The main contribution of this article is presented in Section 3 which contains the algebraic computation. The algebraic techniques are presented: Gröbner basis, a cellular decomposition of semi-algebraic set called cylindrical algebraic decomposition and real roots classification based on Thom's isotopy lemma. They are the computational tools necessary to the algebraic computation in the contrast and multisaturation problems. They are given in details for one specific problem: the classification of the set of singularities of the determinental surface D=0. This shows the complexity of the study (Figs.6 to Figs.11). A final section 4 is devoted to preliminary numerical simulations related to the classification. Non smooth singular arcs are detected in relation with  $D \cap D' = 0$ and preliminary results about the time minimal saturations of two spins are given combining Bocop and HamPath software. Finally LMI techniques implemented via the Globtipoly software are used to evaluate the global optimum in view of validation of the numerical computations of the optimal solutions.

## 2 Geometric frame

The aim of this section is double. First of all, recall the geometric framework relating the classification of singular extremals and the optimal problems. Second, a neat approach is proposed to generate all Lie brackets, in relation with semi-direct Lie algebras.

## 2.1 The classification problem of singular extremals

In this section, we recall the results of [2] which will justify all out algebraic computations.

Definition 1. Let E and F be two  $\mathbb{R}$ -vector spaces, and let G be a group acting linearly on E and F. An homomorphism  $\mathcal{X}: G \to \mathbb{R}^0$  is called a *character*. Let  $\mathcal{X}$  be a character. A *semi-invariant* of weight  $\mathcal{X}$  is a map  $\lambda: E \to \mathbb{R}$  such that for all  $g \in G$  and all  $x \in E$ ,  $\lambda(g.x) = \mathcal{X}(g)\lambda(x)$ . It is an *invariant* if  $\mathcal{X} = 1$ . A map  $\lambda: E \to F$  is a *semi-covariant* of weight  $\mathcal{X}$  is for all  $g \in G$  and for all  $x \in E$ ,  $x \in E$ , x

**Notations and definitions** Let  $U \subset \mathbb{R}^n$  be an open subset and let  $V^\omega(U)$  be the set of real analytic vector fields on U identified to maps from U to U. We denote by  $\mathcal{C}^\omega(U,\mathbb{R})$  the set of analytic maps from U to  $\mathbb{R}$ . Let  $G_d$  be the group of  $C^\omega$  diffeomorphisms of U. The coordinates of  $TU \simeq U \times \mathbb{R}^n$  are denoted z = (q,p) and we endow  $T^*U$  with its canonical symplectif structure defined by  $\Omega = \sum_{i=1}^n \mathrm{d} q_i \wedge \mathrm{d} p_i$ . Let Z be a  $C^\omega$  vector field on U, the Hamiltonian lift of Z is defined by the Hamiltonian  $H_Z(z) = p \cdot Z(q)$ . If  $\varphi \in G_d : q = \varphi(Q)$ , one can lift  $\varphi$  into a symplectic diffeomorphism  $\vec{\varphi}$ , called a *Mathieu transformation* defined by:

$$\vec{\varphi}: q = \varphi(Q), p = P \frac{\partial \varphi^{-1}}{\partial Q}$$

writing p and P as row vectors. A system  $\dot{q} = X + uY$ ,  $X, Y \in V^{\omega}(U)$ , is written as (X, Y) and let  $S = \{(X, Y)\}$ . Take  $(X, Y), (X', Y') \in S$ , they are called *feedback equivalent* if there exists  $\varphi \in G_d$  and a feedback  $u = \alpha(q) + \beta(q)v$ ,  $\alpha, \beta \in C^{\omega}(U, \mathbb{R}), \beta(q)$  non-zero for each q, such that

$$\begin{cases} X' = \varphi * X + \varphi * Y . \alpha \\ Y' = \varphi * Y . \beta \end{cases}$$

where  $\varphi * Z$  is the image of Z defined by  $\varphi * Z = \frac{\partial \varphi^{-1}}{\partial q}(Z \circ \varphi)$ . This action defines a group structure on the set of triplets  $(\varphi, \alpha, \beta)$  and this group is the *feedback group* and denoted  $G_f$ .

Singular extremals of order 2 Let  $(F,G) \in \mathcal{S}$ , the singular extremals of order 2 are the solutions of

$$\dot{z} = \vec{H}_F(z) + u_s \vec{H}_G(z), z \in \Sigma : H_G = 0$$

and such that  $\{\{H_G, H_F\}, H_G\}(z)$  is never vanishing and  $u_s$  is given by (11) as

$$u_s = -\frac{\{\{H_G, H_F\}, H_F\}(z)}{\{\{H_G, H_F\}, H_G\}(z)}$$
(13)

and they are contained in the surface  $\Sigma': H_G(z) = \{H_G, H_F\}(z) = 0$ . Let  $S: \{\{H_G, H_F\}, H_G\} = 0$ . Plugging such  $u_s$  in  $H(z, u) = H_F(z) + uH_G(z)$  defines a true Hamiltonian denoted  $\hat{H}(z)$ . One has the following, see [2].

**Proposition 2.** The singular extremals of order 2 are the solutions of the Hamiltonian equation

$$\dot{q} = \frac{\partial \hat{H}}{\partial p}, \dot{p} = -\frac{\partial \hat{H}}{\partial q}, (q, p) \in \Sigma' \setminus S$$
 (14)

Moreover, an explicit representation is as follows:

**Lemma 1.** Assume that G and [G,F] are never colinear on U. Then  $(\Sigma' \setminus S, \Omega')$  is a symplectif manifold,  $\Omega'$  denoting the restriction of the standard symplectif form and  $\hat{H}$  on  $\Sigma' \setminus S$  is a  $C^{\omega}$  Hamiltonian whose solutions are singular extremals of order 2.

The action on the feedback group  $G_f$  on the set of singular extremals of order 2 is defined as follows: if  $(\varphi, \alpha, \beta) \in G_f$ , then the feedback  $(\alpha, \beta)$  acts trivially and  $\varphi$  acts by the change of symplectic coordinates  $\vec{\varphi}$ .

The key result in our analysis is the following [2].

**Theorem 1.** Let  $\lambda$  be the map  $(F,G) \to (\hat{H},\Sigma',S)$  (differential equation defined by (14)). Then for the respective action of the feedback group  $G_F$ ,  $\lambda$  is a covariant.

Moreover, in relation with our geometric study, it is worth extracting two more covariants, where  $G_F$  acts on functions by  $\vec{\phi}$  action only. This is called a feedback invariant.

**Proposition 3.** The following sets are feedback invariants:

- 1. The set C where F and G are colinear
- 2. The set C' where G and [F,G] are colinear
- 3. On  $\Sigma'$ , the set S defined by  $\{\{H_G, H_F\}, H_G\} = 0$ .

**Reparameterization** In order to study the meromorphic equation (14), one makes a change of parameterization:

$$ds = dt / \{ \{ H_G, H_F \}, H_G \}$$
 (15)

and observes that the map:  $(F,G) \mapsto \{\{H_G,H_F\},H_G\}$  is a *semi*-covariant if  $z \in \Sigma'$ .

**Theorem 2.** The map  $\lambda'$ ;  $(F,G) \mapsto (\hat{H}\{\{H_G,H_F\},H_G\},\Sigma')$  is a semi-covariant.

Note that the associated vector field is  $C^{\omega}$  but not Hamiltonian. It amounts to relate the classification problem to the classification of smooth vector fields and is the basic tool to generate invariants.

**Exceptional and generic cases** The set of singular extremals of order 2 can be split according to [3] into two subsets, recalling that the level sets  $\hat{H} = c$  are invariants.

**Exceptional case** It corresponds to  $\hat{H} = 0$  and the corresponding singular extremals of order 2 are called *exceptional*. Their role in Mayer optimal problems is analyzed in [3].

In our situation, in dimension 4, the constraint  $\hat{H} = 0$  leads to the additional relation  $H_F = p \cdot F = 0$ . Hence the singular control is defined by

$$\begin{cases} H_F = H_G = \{H_G, H_F\} = 0\\ \{\{H_G, H_F\}, H_F\} + u_s\{\{H_G, H_F\}, H_G\} = 0 \end{cases}$$

Since  $p \in \mathbb{R}^4 \setminus \{0\}$  this leads to the relation

$$D'(q) + u_s D(q) = 0$$

where D, D' are the associated determinants

$$D = \det(F, G, [G, F], [[G, F], G])$$
  
$$D' = \det(F, G, [G, F], [[G, F], F])$$

and the corresponding control is given by a feedback  $u_s = -\frac{D'(q)}{D(q)}$  and the vector field is denoted  $X^e(q) = F(q) - \frac{D'(q)}{D(q)}G(q)$ , which again can be analyzed using the reparameterization ds = dt/D(q(t)) and this gives the smooth vector field

$$X_r^e = DF - D'G. (16)$$

In this case the  $\vec{\varphi}$ -action is reduced to the standard action of diffeomorphisms  $\varphi$  on U acting on vector fields V by image  $\varphi * V$  and on mappings  $f: U \to \mathbb{R}$  by composition.

**Generic case** It corresponds to  $\hat{H} = c \neq 0$  and by homogeneity one can take c = 1. Let  $\mathcal{D}' = \{\{H_G, H_F\}, H_G\}, \mathcal{D} = \{\{H_G, H_F\}, H_F\}$ , the singular control  $u_S$  is defined by

$$\begin{cases} H_G = \{H_G, H_F\} = 0 \\ \mathcal{D} + u_S \mathcal{D}' = 0. \end{cases}$$

# 2.2 Lie brackets computations

Each spin system can be lifted on the semi-direct product  $S = GL(3, \mathbb{R}) \times \mathbb{R}^3$  acting on the q-space using the action (A, a).q = Aq + a. The Lie bracket computation rule is [(A, a), (B, b)] = ([A, B], Ab - Ba). Our system is written (F, G) with

- $F = (A, a), A = \operatorname{diag}(-\Gamma, -\Gamma, -\gamma)$  and  $a = (0, 0, \gamma)$
- G = Cq where C is the antisymmetric matrix  $C = E_{32} E_{23}$  with  $E_{ij} = (\delta_{ij})$ .

According to the Lie bracket computation rule on the semidirect Lie product, we can reduce the computations to matrix Lie brackets. Moreover, to make such computations we use the following standard results, see [17] for the details.

The Lie algebra  $\mathfrak{gl}(n,\mathbb{R})$  of  $n \times n$  matrices decomposes into  $cI_n \oplus \mathfrak{sl}(n,\mathbb{R})$  where  $\mathfrak{sl}(n,\mathbb{R})$  is the Lie algebra of matrices with zero trace. This algebra can be written as the direct sum  $a \oplus b$  with a is the Cartan subalgebra of diagonal matrices and  $b = \bigoplus_{i \neq j} \mathbb{R}E_{ij}$ . Moreover we have the Cartan decomposition  $\mathfrak{sl}(n,\mathbb{R}) = \mathfrak{so}(n) \oplus h$  where h is the subspace of symmetric matrices. If  $A = \operatorname{diag}(\lambda_1, \ldots, \lambda_n) \in a$  then  $[A, E_{ij}] = (\lambda_i - \lambda_j)E_{ij}$  and  $\lambda_i - \lambda_j$  will form the nontrivial spectrum of  $\operatorname{ad} A$  (where  $\operatorname{ad} A$  is defined as  $\operatorname{ad} A(B) = [A, B]$ ) with corresponding eigenvectors  $E_{ij}$ .

**Application** Restricting to the q=(y,z) space we have  $A=\operatorname{diag}(-\Gamma,-\gamma)$  and  $a=(0,\gamma)$  and  $A=cI_2+\operatorname{diag}(\lambda,-\lambda)$  where  $c=-\frac{\Gamma+\gamma}{2}$  which is zero if and only if  $\gamma=\Gamma=0$  and  $\lambda=\delta/2$  where  $\delta=\gamma-\Gamma$ . If  $\delta\neq 0$ , the nontrivial spectrum is  $(\delta,-\delta)$ .

In NMR, we have the following

## **Lemma 2.** The case $\gamma = \Gamma$ is the case of water species.

Otherwise, an easy computation gives that the Lie algebra generated by ((A,a),(C,0)) is  $\mathfrak{gl}(2,\mathbb{R})\oplus\mathbb{R}^2$ . Moreover, all Lie brackets can be easily computed. They are listed next, up to order 4, needed in our computation.

## Length 1

- $F = (-\Gamma y, \gamma(1-z))$
- G = (-z, y)

## Length 2

• 
$$[F,G] = (\gamma - \delta z, -\delta y)$$

## Length 3

- $[[F,G],F] = (-\gamma(\gamma-2\Gamma)+\delta^2z,-\delta^2y)$
- $[[F,G],G] = (-2\delta y, -\gamma + 2\delta z)$

## Length 4

- $[[[F,G],F],F] = (-\delta^3 z + \gamma (\gamma^2 3\gamma \Gamma + 3\Gamma^2), -\delta^3 y)$
- $[[[F,G],F],G] = [[[F,G],G],F] = (0,\gamma(\gamma-2\Gamma))$
- $[[[F,G],G],G] = (-\gamma + 4\delta z, 4\delta y)$

## Couple of spins

- Contrast, multisaturation: pairs with  $(\Gamma_1, \gamma_1), (\Gamma_2, \gamma_2)$
- Multisaturation with  $B_1$  inhomogeneity: in this case  $\Gamma_1 = \Gamma_2 = \Gamma$ ,  $\gamma_1 = \gamma_2$  and for the second spin the vector field G is rescaled into  $(1 \varepsilon)G$ .

# 3 Algebraic computations

## 3.1 Preliminaries

**Four test cases** In the physical experiments different cases will be considered. For the contrast problem:

- In vitro four test cases [4] with relaxation times in seconds:
  - $P_1$ : water  $(T_1 = T_2 = 2.5 \text{ s})$  / cerebrospinal fluid  $(T_1 = 2 \text{ s}, T_2 = 0.3 \text{ s})$
  - $P_2$ : deoxygenated blood ( $T_1 = 1.35 \,\text{s}$ ,  $T_2 = 0.05 \,\text{s}$ ) / oxygenated blood ( $T_1 = 1.35 \,\text{s}$ ,  $T_2 = 0.2 \,\text{s}$ )
  - $P_3$ : gray cerebral matter ( $T_1 = 0.92$  s,  $T_2 = 0.1$  s) / white cerebral matter ( $T_1 = 0.780$  s,  $T_2 = 0.09$  s)
  - $P_4$ : water  $(T_1 = T_2 = 2.5 \text{ s})$  / fat  $(T_1 = 0.2 \text{ s}, T_2 = 0.1 \text{ s})$
- *In vivo* one test case [33]:
  - $Q_1$ : brain ( $T_1 = 1.062$  s,  $T_2 = 0.052$  s) / parietal muscle ( $T_1 = 1.200$  s,  $T_2 = 0.029$  s)

Colinear sets (feedback invariants) We denote by  $\Pi_1, \Pi_2$  the respective projections on the first and second spind:  $q \mapsto q_1, q \mapsto q_2$ . Each colinear set is a curve described by the projections corresponding to the colinear set of each spin and an additional relation.

•  $C_1$ : F, G linearly dependent: the projections are the ovals defined by

$$\Gamma_i y_i^2 = \gamma_i (1 - z_i) z_i, 0 \le z_1 \le 1$$

intersected with either of the sets

$$\Gamma_1 y_1 z_2 = \Gamma_2 y_2 z_1$$

$$\gamma_1 y_1 (z_2 - 1) = \gamma_2 y_2 (z_1 - 1)$$

for the contrast setting, or with either of the sets

$$\Gamma(1-\varepsilon)y_1z_2 = \Gamma y_2z_1$$

$$\gamma y_1(z_2-1) = \gamma(1-\varepsilon)y_2(z_1-1)$$

for the  $B_1$ -inhomogeneity setting.

•  $C_2$ : G, [G, F] linearly dependent: for a single spin, the projections are the 2 lines defined by

$$y = 0$$
  $2\delta z = \gamma \text{ (if } \delta \neq 0 \text{, that is if the matter is not water)}$ 

For two spins in the contrast setting, the projections are:

- the curve  $y_1 = y_2 = 0$ ,  $\gamma_1 z_2 \gamma_2 z_1 = (\delta_1 \delta_2) z_1 z_2$ ;
- the 2 lines  $y_i = 0$ ,  $z_1 = \gamma_i/(\delta_1 + \delta_i)$ ,  $z_2 = \gamma_i/(\delta_2 + \delta_i)$  (i = 1, 2);
- the plane  $z_1 = \gamma_1/2\delta$ ,  $z_2 = \gamma_2/2\delta$  if  $\delta_1 = \delta_2 = \delta$ .

For two spins in the  $B_1$ -inhomogeneity setting, the projections are:

- the line  $y_1 = y_2 = 0$ ,  $z_1 = z_2$ ;
- the plane  $z_1 = z_2 = \gamma/2\delta$ .

# 3.2 Algebraic techniques

In this section, we give a brief overview of algebraic tools and methods to solve the algebraic problems posed by the study.

## 3.2.1 Operations on polynomial ideals

Throughout the rest of this section, we let n be a positive integer, and we consider polynomials in  $A = \mathbb{C}[X_1, \dots, X_n]$ .

Our main object of study shall be polynomial ideals. Their importance in algebraic geometry stems from the fact that they are a particularly adequate way to encode a system of equations: given a system of equations F, any system of generators of the ideal  $\langle F \rangle$  has the same zeroes as F, which we will denote V(F) or  $V(\langle F \rangle)$ , and call an algebraic set.

Our work will consist in successive transformations of ideals of polynomial equations, in order to give a human-readable description of them. These transformations are mostly of two forms: elimination of variables, and saturation of polynomials.

Eliminating the first k variables from the ideal I means computing a set of generators of  $I_k := I \cap \mathbb{C}[X_{k+1}, \dots, X_n]$ . Geometrically, the set of zeroes of the ideal  $I_k$  is the smallest algebraic set containing the projection of the set of zeroes V(I) onto the affine subspace  $\mathbb{C}^{n-k}$  with coordinates  $X_{k+1}, \dots, X_n$ .

Given an ideal I and a polynomial f, saturating I by f means computing a set of generators of the ideal

$$(I: f^{\infty}) := \{ g \in A \mid \exists m \in \mathbb{N}, g f^n \in I \}.$$

Geometrically, the zeroes of this ideal is the smallest algebraic set containing  $V(I) \setminus V(f)$ .

This can be computed by adding the polynomial  $u \cdot f - 1$  (with a new variable u) to the generators of I, and eliminating u from that ideal.

This effectively reduces our two major computational tools to the polynomial elimination problem. Because this problem is of such importance in computational algebraic geometry, many tools have been developed for solving it, for example triangular sets [1], resultants [11] or regular chains [27]. In the following computations, we shall mainly use Gröbner bases, because they will allow us to perform both the elimination steps and routine simplification steps.

In the next subsection we give more details about how Gröbner bases can be used in our setting. Before that, we conclude this subsection with a few algebraic and algorithmic techniques which we shall use in the following computations.

Given two systems of polynomial equations F and G, the set  $V(F) \cap V(G)$  is the algebraic set  $V(F \cup G)$ .

The radical of an ideal I is the set

$$\sqrt{I} := \{ f \in A \mid \exists m \in \mathbb{N}, f^m \in I \}.$$

and it has the same set of zeroes as I. Computing the radical of an ideal is a difficult problem, which we will not need to tackle here. However, we shall sometimes be interested in, given a set of generators of I, computing a set of generators of another ideal  $J \supset I$  such that  $\sqrt{I} = \sqrt{J}$ . This can be done by using the square-free form of the generators: given a polynomial f whose (unique) decomposition into primes is  $p_1^{n_1} \cdots p_r^{n_r}$ , its square-free form is  $\mathsf{sqfr}(f) := p_1 \cdots p_r$ .

#### 3.2.2 Gröbner bases

A Gröbner basis of an ideal I is a particular set of generators of I, with additional properties. We only detail the information needed to understand subsequent computations, and we refer the interested reader to [15] for a comprehensive overview of Gröbner basis techniques and elimination theory. The definition of a Gröbner basis depends on the choice of a monomial ordering, that is a total ordering on the monomials  $X_1^{\alpha_1} \cdots X_n^{\alpha_n}$  compatible with multiplication. In the sequel, we will use mostly 3 orderings:

- the *degree reverse-lexicographic* ordering (DRL), denoted  $X_1 > X_2 > \cdots > X_n$ ;
- the *k*'th *elimination ordering* (elim-*k*), denoted  $X_1 > \cdots > X_k \gg X_{k+1} > \cdots > X_n$ , and defined as a block-DRL ordering;
- the *lexicographical* ordering lex, denoted  $X_1 \gg X_2 \gg \cdots \gg X_n$ .

The DRL ordering is usually the easiest to compute, but is nonetheless useful for simple cases, when one merely wants to simplify the generators of an ideal.

The elim-k ordering allows us to eliminate the first k variables: if G is a Gröbner basis of I with respect to the elim-k ordering, then  $G \cap \mathbb{C}[X_{k+1}, \dots, X_n]$  is a (DRL) Gröbner basis of the ideal  $I \cap \mathbb{C}[X_{k+1}, \dots, X_n]$ .

Finally, the lexicographical order has the elimination property for all indices k: from a single Gröbner basis, one can recover all the intermediate elimination bases. This is usually enough to obtain a full description of the solutions, but these bases are also the hardest to compute directly.

Most computer algebra systems include at least one toolbox for Gröbner basis computations. For example in Maple, one can use the built-in package Groebner as well or the third-party library FGb [16].

## 3.2.3 Cylindrical algebraic decomposition

The last two subsections are focussed on two tools for real algebraic geometry. The basic item of study of real algebraic geometry is semi-algebraic sets, that is a sets  $S \subset \mathbb{R}^n$  defined as a finite

union  $S = \bigcup_{i=1}^r S_i$  of sets defined by polynomial equations  $\mathcal{P}_i$  and inequalities  $\mathcal{Q}_i$ :

$$S_i = \left\{ \mathbf{x} \in \mathbb{R}^n \middle| \begin{array}{c} \forall P \in \mathcal{P}_i, P(\mathbf{x}) = 0 \\ \forall Q \in \mathcal{Q}_i, Q(\mathbf{x}) > 0 \end{array} \right\}.$$

Given two semi-algebraic sets A and B, a map  $A \rightarrow B$  is called a semi-algebraic function if its graph is a semi-algebraic set. For example, polynomial or rational functions are semi-algebraic.

A Cylindrical Algebraic Decomposition (CAD) of  $\mathbb{R}^n$  is a sequence  $S_1, \ldots, S_n$  such that for all  $i \in \{1, \ldots, n\}$ ,  $S_i$  is a decomposition of  $\mathbb{R}^i$  into connected semi-algebraic subsets, called cells, and such that:

- Cells of  $S_1$  are all points or intervals;
- For any  $i \in \{1, ..., n\}$ , and for any cell S of  $S_i$ , there exists a finite number of continuous semi-algebraic functions

$$\xi_{S,1},\ldots,\xi_{S,l_S}:S\to\mathbb{R}$$

such that  $S \times \mathbb{R} \subset \mathbb{R}^{i+1}$  is a disjoint union of cells of  $S_{i+1}$ , which are

- either the graph  $\Gamma_{S,j}$  of a function  $\xi_{S,j}$
- or a band  $B_{S,j}$  bounded by  $\xi_{S,j}$  and  $\xi_{S,j+1}$  for  $j \in \{0,\ldots,l_S\}$ , with the convention that  $\xi_{S,0} = -\infty$  and  $\xi_{S,l_S+1} = +\infty$ :

$$B_{S,i} = \{(x', x_{i+1}) \in S \times \mathbb{R} \mid \xi_{S,i}(x') < x_{i+1} < \xi_{S,i+1}(x')\}.$$

Given a set of polynomials  $\mathcal{P} \subset \mathbb{R}[X_1, \dots, X_n]$ , a CAD  $\mathcal{S} = (\mathcal{S}_1, \dots, \mathcal{S}_n)$  is  $\mathcal{P}$ -invariant if, for any cell  $S \in \mathcal{S}_n$  and any polynomial  $P \in \mathcal{P}$ , P has constant sign on S.

In particular, given a semi-algebraic set defined by polynomial equations  $\mathcal{P}$  and polynomial inequalities  $\mathcal{Q}$ , a  $\mathcal{P} \cup \mathcal{Q}$ -invariant CAD gives an extensive topological description of the semi-algebraic set.

Unfortunately, computing a  $\mathcal{P}$ -invariant cylindrical algebraic decomposition, for  $\mathcal{P} \subset \mathbb{Q}[X_1, \dots, X_n]$ , is intrinsically difficult: the complexity of the algorithm and the size of the output, for most entries, is doubly exponential in n.

Nonetheless, this tool will prove useful for completing computations reduced to a smallerdimensional set, for example from the algorithms in the next section. In this context, the main purpose of the CAD will be to compute one point per connected component of an open semialgebraic set, defined by inequations  $P_i \neq 0$  and strict inequalities  $P_i > 0$ .

Cylindrical algebraic decompositions are implemented in Maple as part of the subpackage SemialgebraicsetTools of RegularChains. Implementations also exist for more specific problems. In particular for computing one point per connected component of an open semi-algebraic set, we use the third-party Maple library RAGLib.

## 3.2.4 Real roots classification

Our last algebraic tool is a computational technique for solving the problem of classifying the real fibers of projection restricted to an algebraic set.

Consider an algebraic set V in  $\mathbb{C}^{n+t}$ , with coordinates  $(\mathbf{X}, \mathbf{G}) = (X_1, \dots, X_n, G_1, \dots, G_t)$  and the projection onto the parameter space

$$\pi : \mathbb{C}^{n+t} \rightarrow \mathbb{C}^t$$
 $(\mathbf{x}, \mathbf{g}) \mapsto \mathbf{g}$ 

Assume that V has dimension t. Let  $B \in \mathbb{R}^{n+t}$  be a semi-algebraic set defined only by a set of polynomial inequalities H:

$$B = \{(\mathbf{x}, \mathbf{g}) \in \mathbb{R}^{n+t} \mid \forall h \in H, h(\mathbf{x}, \mathbf{g}) > 0\},\$$

we consider the semi-algebraic set  $V \cap B \subset \mathbb{R}^{n+t}$ .

The classification problem is to identify a dense covering of the *real* parameter space  $\mathbb{R}^t$  into open subsets  $U_1, \ldots, U_N$  such that on each  $U_i$ , the number of points in the fibers of  $\pi$  restricted to  $V \cap B$  is constant:

$$\exists c_i, \forall \mathbf{g} \in U_i, \#\pi^{-1}(\mathbf{g}) \cap V \cap B = c_i.$$

This problem has been thoroughly studied, see [36, 26]. In our situation, the variety V shall be defined as the zero set of D and its partial derivatives along the  $X_i$ 's, where D is the determinant of a matrix. The strategy that we use, as described in [9], is adapted from the more general strategy of described in [36, 26].

We summarize the key points of this general strategy. One wishes to compute a set of polynomials  $\{p_1, \ldots, p_M\} \subset \mathbb{R}[G]$  such that the dense covering  $(U_i)$  given by the connected components of  $\mathbb{R}^t \setminus (V(p_1) \cup V(p_2) \cup \cdots \cup V(p_M))$ . These polynomials correspond to equations satisfied by points at which the cardinality of the fibers change. These points may be:

- the projection of singular points of V;
- the critical values of  $\pi$  restricted to V;
- the projection of points where V meets the topological border  $\partial B$  of the semi-algebraic set B:
- points **g** at which the fiber  $\pi^{-1}(\mathbf{g}) \cap V$  has positive real dimension;
- points at which the projection  $\pi$  restricted to  $V \cap B$  is not proper.

This is formalized in the following lemma, whose proof relies on Thom's isotopy lemma. We first set some notations and hypotheses. If B is a semi-algebraic set defined by polynomial inequalities H, we define

$$B_0 := \bigcup_{h \in H} V(h).$$

Note in particular that  $\partial B \subset B_0$ . We denote by  $\operatorname{sing}(V)$  the singular locus of V,  $\operatorname{crit}(\pi, V)$  the critical locus of  $\pi$  restricted to V, and  $K(\pi, V) = \operatorname{sing}(V) \cup \operatorname{crit}(\pi, V)$ .

The hypotheses are:

 $\mathcal{H}1$  There exists a nonempty Zariski-open subset  $\mathcal{O}_1$  of  $\mathbb{C}^t$  such that for all  $\mathbf{g} \in \mathcal{O}_1$ , the fiber  $V \cap \pi^{-1}(\mathbf{g})$  is a nonempty finite subset of  $\mathbb{C}^{n+t}$ 

- $\mathcal{H}2$  The restriction of the projection  $\pi$  to B is proper;
- $\mathcal{H}3$  The intersection  $V \cap B_0$  has dimension at most t-1 in  $\mathbb{C}^{nt}$ ;
- $\mathcal{H}4$  The algebraic set V is equidimensional with dimension t.

**Lemma 3.** Let V and B satisfying hypotheses  $\mathcal{H}1$ ,  $\mathcal{H}2$ ,  $\mathcal{H}3$  and  $\mathcal{H}4$ . Let U be a nonempty connected open subset of  $\mathbb{R}^t$  which does not meet  $\pi(V \cap B_0) \cup \pi(K(\pi, V))$ , and let  $\mathbf{g} \in U$ . Then  $V \cap \pi^{-1}(\mathbf{g})$  is finite, and for any  $\mathbf{g}' \in U$ ,  $\#V \cap \pi^{-1}(\mathbf{g}') = \#V \cap \pi^{-1}(\mathbf{g})$ .

Hence, under these hypotheses, it suffices to compute a nonzero polynomial whose zeroes cover  $\pi(V \cap B_0)$  and  $\pi(K(\pi, V))$  to obtain the dense covering that we need.

The strategy that we use refines these computations by taking advantage of the fact that our variety V comes from a determinantal variety, defined as follows. Let M be a  $k \times k$  matrix with coefficients in  $\mathbb{R}[\mathbf{X}, \mathbf{G}]$ ,  $r \in \{0, \dots, k-1\}$ , the r'th determinantal variety associated with M is

$$V_r(M) = \{(\mathbf{x}, \mathbf{g}) \in \mathbb{C}^{n+t} \mid \operatorname{rank}(M(\mathbf{x}, \mathbf{g})) \le r\}.$$

These varieties can be defined algebraically in two ways. First, one can use the minors modelization:

$$V_r(M) = V((r+1)$$
-minors of  $M$ ).

Another way to define  $V_r$  is as the projection of the incidence variety defined as the set of solutions of

$$M(\mathbf{x}, \mathbf{g}) \cdot Y(\mathbf{y}) = 0$$

where  $Y(\mathbf{y})$  is a full-rank  $k \times k - r$  matrix whose entries are new variables  $\mathbf{y} = y_{1,1}, \dots, y_{k,k-r}$ . We now describe more precisely how this property allows us to compute the classification for the MRI problem. The variety V that we will consider is defined as  $K(\pi, V_{r_0}(M))$  for  $r_0 \in \{0, \dots, k-1\}$ . Assume that the variety V satisfies the following three properties:

 $\mathcal{H}5$  There exists a nonempty Zariski-open subset  $\mathcal{O}_2 \subset \mathbb{C}^t$  such that

$$V \cap \pi^{-1}(\mathcal{O}_2) = V_{r_0-1} \cap \pi^{-1}(\mathcal{O}_2);$$

- $\mathcal{H}6$  For any  $r_1 < r_2 \in \{0, \dots, k-1\}$ , for any matrix A of size  $r_1$  in M, the ideal generated by the  $r_2$ -minors of M containing A is radical;
- $\mathcal{H}7$  For any  $r \in \{0, \dots, k-1\}$ , the variety  $V_r$  is equidimensional with dimension  $n+t-(k-r+1)^2$ .

These properties are generic [10, 32]. If we assume that

$$n = (k - r_0 + 1)^2,$$

then hypotheses  $\mathcal{H}7$  and  $\mathcal{H}5$  imply hypotheses  $\mathcal{H}4$  and  $\mathcal{H}1$  for V. Under these hypotheses, the strategy consists in computing:

- 1. the projection of  $V \setminus V_{r_0-1}$ , which has dimension less than t;
- 2. the projection of  $K(\pi, V_{r_0-1})$ ;
- 3. the projection of  $B_0 \cap V_{r_0-1}$ .

*Remark* 1. In section 3.3.3, we will also be describing  $K(\pi, V)$  where V is an incidence variety defined by the equations

$$P^t \cdot M(\mathbf{x}, \mathbf{g}) = 0$$

where P is a vector of dimension 4 and M is a matrix of size  $4 \times 3$ . This incidence variety encodes an underdefined linear system, so its projection onto the affine space  $\mathbb{C}^{n+t}$  with coordinates  $(\mathbf{X}, \mathbf{G})$  is the affine space itself. In this situation, generically,  $\operatorname{sing}(V)$  is exactly  $V_2(M)$ .

# 3.3 Algebraic computations for multisaturation with $B_1$ -inhomogeneity

#### 3.3.1 Conventions and notations

In this section and the next one, we use the change of coordinates

$$z_1 \leftarrow z_1 - 1, z_2 \leftarrow z_2 - 1,$$

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).

The following polynomials will appear frequently in the remainder of the section:

- $P_{y_1} := y_1 (1 \varepsilon)y_2$
- $P_{y_2} := y_2 (1 \varepsilon)y_1$
- $P_{z_1} := 2(\Gamma \gamma)z_1 + 2\Gamma \gamma$
- $P_{z_2} := 2(\Gamma \gamma)z_2 + 2\Gamma \gamma$

The root of the univariate polynomials  $P_{z_1}$  and  $P_{z_2}$  is

$$z_S = \frac{\gamma - 2\Gamma}{2\Gamma - 2\gamma}.$$

## 3.3.2 Transfer time not fixed $(H_F = 0)$

Singularities of  $\{D=0\}$ 

**Proposition 4.** The set of points satisfying  $D = \frac{\partial D}{\partial y_1} = \frac{\partial D}{\partial z_1} = \frac{\partial D}{\partial y_2} = \frac{\partial D}{\partial z_2} = 0$  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

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

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

*Proof.* The determinant D can be factored as  $(1 - \varepsilon)\tilde{D}$ . The singularities of D and those of  $\tilde{D}$  are the same, so for the study, we consider the ideal

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

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 order  $y_1 > y_2 > z_1 \gg z_2 > \varepsilon > \Gamma > \gamma$ . This computation yields that

$$I \cap \mathbb{Q}[z_2, \varepsilon, \Gamma, \gamma] = \langle \varepsilon^2(\varepsilon - 2)^2(2\Gamma - \gamma)(\Gamma - \gamma)z_2^3 P_{z_2}^3 \rangle,$$

so singular points necessarily satisfy

$$\begin{cases} z_2 = 0 \\ \text{or} \\ z_2 = \frac{\gamma - 2\Gamma}{2\Gamma - 2\gamma}. \end{cases}$$

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  $2\Gamma \ge \gamma$ ), 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

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

In order to lift the partial solution  $z_2 = 0$ , we compute a Gröbner basis  $G_1$  of  $I_1$  with respect to the order  $y_1 > y_2 \gg z_1 > z_2 > \varepsilon > \Gamma > \gamma$ , and we find that this ideal contains

$$\gamma z_1^2 (\varepsilon - 1)^2 (2\Gamma - \gamma),$$

so  $z_1 = 0$ .

We then compute a Gröbner basis of  $\langle \mathsf{sqfr}(G_1), z_1 \rangle$  with respect to the order  $y_1 \gg y_2 > z_1 > z_2 > \varepsilon > \Gamma > \gamma$ , and we find that this ideal contains

$$\Gamma \gamma \varepsilon y_2^2 (\Gamma - \gamma) (2\Gamma - \gamma)^2 (\varepsilon - 2),$$

so  $y_2 = 1$ .

Finally, adding  $y_2$  to the ideal yields that

$$0 = \Gamma y_1(\varepsilon - 1)(2\Gamma - \gamma),$$

so the final 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  $y_1 > y_2 \gg z_1 > z_2 > \varepsilon > \Gamma > \gamma$ , 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  $\mathsf{sqfr}(G_2)$  and computing a Gröbner basis for the order  $y_1 \gg y_2 > z_1 > z_2 > \varepsilon > \Gamma > \gamma$ , we find that the ideal contains

$$\gamma^2 y_2(\varepsilon-1)P_{y_1}(2\Gamma-\gamma)$$
,

so we have 2 new branches to consider.

If  $y_2 \neq 0$ ,  $y_1 = (1 - \varepsilon)y_2$ . 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  $\{P_{y_1} = 0\}$ .

Locus of  $\{D = D' = 0\}$ 

**Proposition 5.** The points of  $\{D = D' = 0\}$  are given by:

1. the plane

$$z_1 = z_2 = z_S = \frac{\gamma - 2\Gamma}{2\Gamma - 2\gamma} \tag{17}$$

2. the line

$$\begin{cases} y_1 = y_2 = 0 \\ z_1 = z_2 \end{cases}$$
 (18)

3. the surface (parameterized by  $y_1, y_2$ )

$$z_1 = z_2 = \frac{\Gamma P_{y_2}^2 (\gamma - 2\Gamma)}{2(\Gamma - \gamma)a_3}$$
 (19)

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with

$$a_3 = (\Gamma + \gamma)P_{y_1}^2 + \varepsilon(\varepsilon - 2)\Gamma(y_1 - y_2)(y_1 + y_2)$$

4. the surface (parameterized by  $y_1, z_2$ )

$$y_2 = \frac{y_1 z_2}{(1 - \varepsilon)z_1}, \ z_1 = \frac{(2\Gamma - \gamma)z_2}{a_4}$$
 (20)

with

$$a_4 = 2(\varepsilon - 2)(\Gamma - \gamma)\varepsilon z_2 + (2\Gamma - \gamma)(\varepsilon - 1)^2$$

5. the surface (parameterized by  $y_2, z_2$ )

$$z_{1} = \frac{z_{2}y_{1}}{(1 - \varepsilon)y_{2}}, y_{1} = \frac{(1 - \varepsilon)y_{2}\left((2\Gamma - \gamma)\Gamma y_{2}^{2} + \gamma^{2}z_{2}^{2}\right)}{a_{5}}$$
(21)

with

$$a_5 = \Gamma \left( 2\varepsilon (\varepsilon - 2)(\Gamma - \gamma)z_2 + (\varepsilon - 1)^2 (2\Gamma - \gamma) \right) y_2^2 + \gamma^2 z_2^2$$

*Proof.* The determinant D' factors as

$$D' = 2\gamma^2 (2\Gamma - \gamma)(\Gamma - \gamma)(z_1 - z_2)(\varepsilon - 1)((\varepsilon - 1)y_2z_1 + y_1z_2),$$

so we form the two ideals

$$I_1 = \langle \tilde{D}, z_1 - z_2 \rangle$$
  

$$I_2 = \langle \tilde{D}, (\varepsilon - 1)y_2z_1 + y_1z_2 \rangle.$$

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

$$2(\Gamma - \gamma) \left( (\Gamma + \gamma)P_{y_1}^2 + \varepsilon(\varepsilon - 2)\Gamma(y_1 - y_2)(y_1 + y_2) \right) z_2 + \Gamma P_{y_2}^2 (2\Gamma - \gamma)$$

The polynomial  $P_{z_2}$  gives solution 1.

Let

$$a_3(y_1, y_2) = (\Gamma + \gamma)P_{y_1}^2 + \varepsilon(\varepsilon - 2)\Gamma(y_1 - y_2)(y_1 + y_2)$$

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

$$-4(\varepsilon-2)^2\varepsilon^2y_1^2\gamma\Gamma$$
.

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  $(\varepsilon - 1)y_2z_1 + y_1z_2 = 0$ .

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  $2\Gamma \ge \gamma$ , the only solution is  $z_2 = 0$  with either  $2\Gamma = \gamma$  or  $y_2 = 0$ .

If  $y_1 = 0$  and  $z_1 \neq 0$ , then (since  $\varepsilon \neq 1$ ) 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

$$0 = p_3 := (2(\varepsilon - 2)(\Gamma - \gamma)\varepsilon z_2 + (2\Gamma - \gamma)(\varepsilon - 1)^2)z_1 - (2\Gamma - \gamma)z_2.$$

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

$$\begin{cases} a_4(z_2) \neq 0 \\ z_1 = \frac{2\Gamma - \gamma}{c_3(z_2)} z_2 \end{cases}$$

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  $u \gg z_1 > y_1 \gg z_2 > y_2 > \varepsilon > \gamma > \Gamma$ . This basis contains a polynomial which factors as the product of

$$p_4 = (2(\varepsilon - 2)(\Gamma - \gamma)\varepsilon z_2 + (\varepsilon - 1)^2(2\Gamma - \gamma))y_1 + (\varepsilon - 1)(2\Gamma - \gamma)y_2$$

and

$$p_5 = \left(\Gamma\left(2\varepsilon(\varepsilon-2)(\Gamma-\gamma)z_2 + (\varepsilon-1)^2(2\Gamma-\gamma)\right)y_2^2 + \gamma^2z_2^2\right)y_1 + (\varepsilon-1)y_2\left((2\Gamma-\gamma)\Gamma y_2^2 + \gamma^2z_2^2\right).$$

First, assume that  $p_4 = 0$ . We compute a Gröbner basis of  $I_2 + \langle uy_1 - 1, p_4 \rangle$  for the order  $u \gg z_1 > y_1 \gg z_2 > y_2 > \varepsilon > \gamma > \Gamma$ , 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

$$-4a_4(z_2)\Gamma\gamma^2z_2^2$$

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

$$\begin{cases} a_5(y_2, z_2) \neq 0 \\ y_1 = \frac{(1-\varepsilon)y_2((2\Gamma-\gamma)\Gamma y_2^2 + \gamma^2 z_2^2)}{\Gamma(2\varepsilon(\varepsilon-2)(\Gamma-\gamma)z_2 + (\varepsilon-1)^2(2\Gamma-\gamma))y_2^2 + \gamma^2 z_2^2} \end{cases}$$

and

$$\begin{cases} a_5(y_2, z_2) = 0 \\ y_2((2\Gamma - \gamma)\Gamma y_2^2 + \gamma^2 z_2^2) \end{cases}$$

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.

## Equilibrium positions

**Lemma 4.** 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.

**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

$$\frac{\mathrm{d}}{\mathrm{d}t}(q+\delta q) = (DF - D'G)(q) + A(q) \cdot \delta q + R(\delta q).$$

where  $A = \operatorname{Jac}_q(DF - D'G)$ , so that

$$\frac{\mathrm{d}}{\mathrm{d}t}(\delta q) = A(q) \cdot \delta q + R(q)(\delta q).$$

We can compute A(q) explicitely: Indeed, let f = DF - D'G. Its first derivative is

$$df(q)(u) = dD(q)(u)F(q) + D(q)dF(q)(u) - dD'(q)(u)G(q) - D'(q)dG(q)(u),$$
(22)

so

$$A(q) = \nabla D(q).F(q) + D(q)\operatorname{Jac}_{q}(F)(q) - \nabla D'(q).G(q) - D'(q)\operatorname{Jac}_{q}(G)(q).$$

We examine the eigenvalue decomposition of A(q).

**Solution 1** (17) If  $z_1 = z_2 = \frac{\gamma - 2\Gamma}{2\Gamma - 2\gamma}$ , the characteristic polynomial of A factors as

$$T^2 \left(T - \gamma^2 (2\Gamma - \gamma)^2 (\varepsilon - 1) P_{y_1}^2\right)^2$$

The matrix A(q) is diagonalizable.

**Solution 2** (18) If  $y_1 = y_2 = 0$  and  $z_1 = z_2$ , the characteristic polynomial of A(q) is

The Jacobian matrix A(q) can be trigonalized as

with the transition matrix

$$P = \begin{bmatrix} 0 & 1 & -1 & 0 \\ \varepsilon \gamma^3 (\varepsilon - 1)(\varepsilon - 2) z_1^2 P_{z_1} & 1 & 0 & 0 \\ 0 & 1 & 0 & 1 \\ \varepsilon \gamma^3 (\varepsilon - 1)(\varepsilon - 2) z_1^2 P_{z_1} & 0 & 0 & 0 \end{bmatrix}$$

**Solution 3** (19) If  $z_1 = z_2 = \Gamma P_{y_2}^2 (\gamma - 2\Gamma)/2 (\Gamma - \gamma) a_3$ , 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

$$b_3 = \Gamma \gamma^2 (\varepsilon - 1) P_{y_1} P_{y_2} (2\Gamma - \gamma)^2.$$

The matrix A(q) is diagonalizable.

**Solution 4** (20) If  $y_2 = \frac{y_1 z_2}{(1-\varepsilon)z_1}$  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

$$b_4 = 2\varepsilon^2 \gamma^3 z_2^3 (\varepsilon - 1)(\varepsilon - 2)^2 (2\Gamma - \gamma)(\Gamma - \gamma) P_{z_2}.$$

The matrix A(q) is diagonalizable.

**Solution 5** (21) If  $z_1 = \frac{z_2y_1}{(1-\varepsilon)y_2}$  and  $y_1 = \frac{(1-\varepsilon)y_2\left((2\Gamma-\gamma)\Gamma y_2^2+\gamma^2 z_2^2\right)}{a_5}$ , 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:

$$\frac{\mathrm{d}}{\mathrm{d}t}(N+\delta q) = R(N)(\delta q) = O(\|\delta q\|^3).$$

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):

$$\frac{\mathrm{d}}{\mathrm{d}t}(q+\delta q) = (DF - D'G)(q) + A(q)(\delta q) + Q(q)(\delta q) + O(\|\delta q\|^3),$$

with  $\frac{dq}{dt}(S) = (DF - D'G)(S) = 0$  and A(S) = 0. We can compute Q by differentiating f = DF - D'G again, as was done in [4, Sec. 3.4]. Differentiating (22) along q again, the second derivative of f is

$$d^{2}f(q)(u,v) = d^{2}D(q)(u,v)F(q) + dD(q)(u)dF(q)(v) + dD(q)(v)dF(q)(u) - d^{2}D'(q)(u,v)G(q) - dD'(q)(u)dG(q)(v) - dD'(q)(v)dG(q)(u)$$
(23)

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

We wish to compute  $H_2(\delta q) = Q(S)(\delta q, \delta q) = \frac{1}{2}d^2f(q)(\delta q, \delta q)$ . 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

$$\begin{split} F(S) &= \left(0, \frac{\gamma(2\Gamma - \gamma)}{2(\Gamma - \gamma)}, 0, \frac{\gamma(2\Gamma - \gamma)}{2(\Gamma - \gamma)}\right)^t \\ G(S) &= \left(\frac{\gamma}{2(\Gamma - \gamma)}, 0, \frac{(1 - \varepsilon)\gamma}{2(\Gamma - \gamma)}, 0\right)^t \\ h_2(\delta q) &= \frac{1}{2} \mathrm{d}^2 D(S)(\delta q, \delta q) = (1 - \varepsilon)(\delta z_1 - \delta z_2)(\delta z_1 - (1 - \varepsilon)^2 \delta z_2)(2\Gamma - \gamma)\gamma^2 \\ h_2'(\delta q) &= \frac{1}{2} \mathrm{d}^2 D'(S)(\delta q, \delta q) = (1 - \varepsilon)(\delta z_1 - \delta z_2)(\delta y_2(\varepsilon - 1) + \delta y_1)(2\Gamma - \gamma)^2 \gamma^2. \end{split}$$

Following [4] and [28], 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

$$\dot{w} = \frac{1}{\|v\|^2} \left( \dot{v} \|v\| - v \frac{\langle v, \dot{v} \rangle}{\|v\|} \right)$$

$$= \frac{H_2(v)}{\|v\|} - \frac{\langle v, H_2(v) \rangle}{\|v\|^3} v$$

$$= \|v\| \left( H_2(w) - \langle w, H_2(w) \rangle w \right)$$

so we are to study the following differential equation on the sphere  $S^3$ :

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

Invariants are related to the eigenvalues of the linearization of  $H_2^{\pi}$  at points where  $H_2^{\pi}(v) = 0$ . 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  $H_2^{\pi}$  in some neighborhood of these solutions in  $S^3$ .

**Proposition 6.** The blow-up at point S has no ray solutions, and two sets of non-isolated singularities:

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

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

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

*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  $H_2(\delta q) = 0$ , and 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 study the linearization of  $H_2^{\pi}$  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  $\bar{\delta y_1} = \delta y_1/\delta z_1$ ,  $\bar{\delta y_2} = \delta y_2/\delta z_1$ ,  $\bar{\delta z_2} = \delta z_2/\delta z_1$ . The differential equation becomes

$$\frac{\mathrm{d}}{\mathrm{d}t} \begin{pmatrix} \delta \bar{y}_1 \\ \delta \bar{y}_2 \\ \delta \bar{z}_1 \end{pmatrix} = \delta z_2 \bar{C}(\delta \bar{y}_1, \delta \bar{y}_2, \delta \bar{z}_1)$$

with  $\bar{C}$  a polynomial vector field of degree 3.

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

$$\begin{bmatrix} 0 & 0 & (\varepsilon-1)(\bar{\delta y_1}(\varepsilon(\varepsilon-1)-1)+\bar{\delta y_2}(1-\varepsilon))(2\Gamma-\gamma)^2\gamma^3 \\ 0 & 0 & (\varepsilon-1)(\bar{\delta y_1}(2\varepsilon(\varepsilon-2)+1)+\bar{\delta y_1}(\varepsilon-1))(2\Gamma-\gamma)^2\gamma^3 \\ 0 & 0 & 0 \end{bmatrix}$$

Then we consider vectors  $\delta q$  such that  $\delta y_2 = (1 - \varepsilon)\delta y_1$ ,  $\delta z_1 = (1 - \varepsilon)^2 \delta z_2$ . This time we use the chart  $\delta z_2 \neq 0$  with coordinates  $\delta \bar{y}_1 = \delta y_1/\delta z_2$ ,  $\delta \bar{y}_2 = \delta y_2/\delta z_2$ ,  $\delta \bar{z}_1 = \delta z_1/\delta z_2$ . As above, we compute the differential equation in this chart, linearize the resulting vector field, and evaluate this Jacobian at  $\delta \bar{y}_2 = (1 - \varepsilon)\delta \bar{y}_1$  and  $\delta \bar{z}_1 = (1 - \varepsilon)^2$ . This matrix has rank 2 and is diagonalizable with non-zero eigenvalues:

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

**Cubic approximation at** N We perform the same study at the North pole N. With expression (23), 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 7.** 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

$$\begin{cases} \delta z_1 = \delta z_2 = 0\\ (\varepsilon - 1)\delta y_1 + \delta y_2 = \frac{1}{\Gamma(2\Gamma - \gamma)\sqrt{1 - \varepsilon}}; \end{cases}$$
 (24)

2. the quadric

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

and three sets of real non-isolated singularities:

1. the plane

$$\begin{cases} \delta z_1 = \delta z_2 \\ \delta y_1 (1 - \varepsilon) = \delta y_2 \end{cases}$$
 (25)

2. the plane

$$\begin{cases} \delta y_2 = (1 - \varepsilon) \delta y_1 \\ \delta z_2 = (1 - \varepsilon)^2 \delta z_1 \end{cases}$$
 (26)

3. the surface defined by

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

$$(27)$$

For points on the line (24), the linearization of  $H_3^{\pi}$  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 (25), 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 (26), 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

$$\left(\Gamma^2(\varepsilon-1)(2\Gamma-\gamma)^2\left((\varepsilon-1)\delta y_1+\delta y_2\right)^2\right)^2-1=0.$$

Since  $\varepsilon - 1 < 0$ ,  $2\Gamma - \gamma > 0$  and  $\Gamma > 0$ , this defines 2 lines of real solutions given by

$$(\varepsilon - 1)\delta y_1 + \delta y_2 = \pm \frac{1}{\Gamma(2\Gamma - \gamma)\sqrt{1 - \varepsilon}}$$

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

$$\left(\gamma^3(2\Gamma-\gamma)(\varepsilon-1)\left(((\varepsilon-1)\delta z_1-\delta z_2)^2-(\varepsilon-2)\delta z_1\delta z_2\right)\right)^2-1=0,$$

which defines 1 quadric

$$((\varepsilon-1)\delta z_1 - \delta z_2)^2 - (\varepsilon-2)\delta z_1\delta z_2 = \frac{1}{\gamma^3(2\Gamma-\gamma)(1-\varepsilon)}.$$

We now consider non-isolated singularities, that is zeroes of  $H_3$ . To this end, we compute a Gröbner basis of the system  $H_3(\delta q) = 0$ , saturating by  $\gamma - \Gamma$ ,  $\gamma$ ,  $\Gamma$ ,  $\varepsilon - 1$  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.  $\delta z_2 = (1 \varepsilon)^2 \delta z_1$
- 5. otherwise.

In the case 1, a Gröbner basis is given by

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

and the solutions form the plane (25).

In the case 2, saturating by  $\delta z_1 - \delta z_2$ , a Gröbner basis is given by

$$\begin{cases} \delta y_1 + (1-\varepsilon)\delta y_2 \\ \delta z_1 + \delta z_2 \\ \Gamma(\varepsilon^2 - 2\varepsilon + 2)(2\gamma - \gamma)\delta y_2^2 + 2\gamma^2\delta z_2^2 \end{cases}$$

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

$$\{\delta y_1\delta z_2 + (\varepsilon - 1)\delta y_2\delta z_1, \delta y_2 + (\varepsilon - 1)\delta y_1, \delta z_2 - (\varepsilon - 1)^2\delta z_1\}$$

and the solutions form the plane (26).

Finally, for the case 5, we compute a Gröbner basis, saturating by all the previous conditions. This basis is

$$\left\{ \begin{aligned} &\Gamma(\varepsilon-1)(2\Gamma-\gamma)\delta y_1^2 + (2\Gamma-\gamma)\Gamma\delta y_1\delta y_2 + \gamma^2(\varepsilon-1)\delta z_1^2 - \gamma^2(\varepsilon-1)\delta z_1\delta z_2 \\ &\Gamma(\varepsilon-1)(2\Gamma-\gamma)\delta y_1\delta y_2 - \gamma^2\delta z_1\delta z_2 + (2\Gamma-\gamma)\Gamma\delta y_2^2 + \gamma^2\delta z_2^2 \\ &\delta y_1\delta z_2 + (\varepsilon-1)\delta y_2\delta z_1 \\ &\left(\Gamma(2\Gamma-\gamma)(\varepsilon-1)^2\delta y_2^2 + \delta z_2^2\gamma^2\right)\delta z_1 - (2\Gamma-\gamma)\Gamma\delta z_2\delta y_2^2 - \gamma^2\delta z_2^3 \end{aligned} \right\}$$

The fourth polynomial is a combination of the other 3, and the solutions form the surface (27).

For the second part of the proposition, as in the quadratic case, we study the linearization of  $H_3^{\pi}$ :  $\dot{v} = H_3(v) - \langle v, H_3(v) \rangle v$ . In the affine chart with  $\delta y_1 \neq 0$ , with coordinates  $\delta \bar{z}_1 = \delta z_1/\delta y_1$ ,  $\delta \bar{y}_2 = \delta y_2/\delta y_1$  and  $\delta \bar{z}_2 = \delta z_2/\delta y_1$ , the differential equation  $\dot{v} = H_3(v)$  becomes

$$\frac{\mathrm{d}}{\mathrm{d}t} \begin{pmatrix} \bar{\delta z_1} \\ \bar{\delta y_2} \\ \bar{\delta z_2} \end{pmatrix} = \delta y_1^2 \bar{Q}(\bar{\delta z_1}, \bar{\delta y_2}, \bar{\delta z_2})$$

with  $\bar{Q}$  a polynomial vector field of degree 4. In this chart,  $H_3^{\pi}$  becomes

$$\dot{v} = \bar{Q}(v)$$
.

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

## 3.3.3 General case

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

**Proposition 8.** The set of singularities of  $\{\mathcal{D} = H_G = \{H_G, H_F\} = 0\}$  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  $z_1 = z_2$ ,  $y_2 = (1 - \varepsilon)y_1$ .

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

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

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

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

has rank at most 2 (as per Remark 1).

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  $\mathbf{L} = L_1, L_2, L_3$  and random numbers  $v_1, v_2, v_3$ . This gives us a system of 4 polynomial equations in the 10 unknowns  $y_1, z_1, y_2, z_2, L_1, L_2, L_3, \Gamma, \gamma, \varepsilon$ . We eliminate  $L_1, L_2, L_3$  from the ideal in order to recover the projection, and we saturate by  $1 - \varepsilon$ .

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

$$(z_1-z_2)((\varepsilon-1)P_{z_2}y_1+P_{z_1}y_2).$$

First, we add  $z_1 - z_2$  to the ideal. Once again, we compute a Gröbner basis for the elimination order  $y_1 > y_2 > z_1 > z_2 \gg \gamma > \Gamma > \varepsilon$ , 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  $(\varepsilon - 1)y_1 + y_2$  to the ideal and saturating by  $P_{z_2}$ ,  $\varepsilon$  and  $\varepsilon - 2$ )

$$\begin{cases} 0 &= y_1(\Gamma - \gamma) \\ 0 &= y_2(\Gamma - \gamma) \\ 0 &= P_{y_2} = (\varepsilon - 1)y_1 + y_2 \\ 0 &= z_1 - z_2. \end{cases}$$

The latter is generically (if  $\Gamma \neq \gamma$ ) defined by

$$\begin{cases} y_1 = y_2 = 0 \\ z_1 = z_2 \end{cases}$$

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

$$\begin{cases} y_1 = \frac{y_2}{\varepsilon - 1} \\ z_1 = z_2. \end{cases}$$

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

Locus of  $\{D = D' = H_G = \{H_G, H_F\} = 0\}$ 

**Proposition 9.** The solutions form the union of the hyperplane defined by

$$z_1 = z_2 \tag{28}$$

and the hypersurface

$$y_1 = -\frac{y_2 P_{z_1}}{(\varepsilon - 1) P_{z_2}} \tag{29}$$

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

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

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

$$\begin{split} \Delta' &= \det \begin{bmatrix} G & [G,F] & [[G,F],G] & [[G,F],F] \\ (1-\varepsilon)G & [(1-\varepsilon)G,F] & [[(1-\varepsilon)G,F],(1-\varepsilon)G & [[(1-\varepsilon)G,F],F] \end{bmatrix} \\ &= (1-\varepsilon)^2 \det \begin{bmatrix} G & [G,F] & [[G,F],G] & [[G,F],F] \\ G & [G,F] & (1-\varepsilon)[[G,F],G & [[G,F],F]. \end{bmatrix} \end{split}$$

This determinant factors as

$$-2(\varepsilon-1)^{2}(2\Gamma-1)(\Gamma-1)(z_{1}-z_{2})[(\varepsilon-1)P_{z_{1}}y_{1}+P_{z_{1}}y_{2}]$$

## Equilibrium positions

**Lemma 5.** All equilibrium points of  $\dot{Z} = \mathcal{D}\vec{H}_F - \mathcal{D}'\vec{H}_G$  satisfying  $H_G = \{H_G, H_F\} = 0$  are contained in  $\{\mathcal{D} = \mathcal{D}' = 0\}$ .

*Proof.* Recall that  $\vec{H}_F$  is defined as

$$ec{H_F} = egin{bmatrix} rac{\partial H_F}{\partial p} \ -rac{\partial F}{\partial q} \end{bmatrix} = egin{bmatrix} F \ -p \cdot rac{\partial H_F}{\partial q} \end{bmatrix}$$

and  $\vec{H}_G$  is defined in the same way. Let z = (q, p) be a point such that

$$\mathcal{D}(z)\vec{H}_F(z) - \mathcal{D}'(z)\vec{H}_G(z) = 0,$$

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  $X_{\mathcal{D}}$  and  $X_{\mathcal{D}'}$  and consider the ideal generated by

- $X_{\mathcal{D}}F(q) X_{\mathcal{D}'}G(q) = 0$
- $H_G(z) = \{H_G, H_F\}(z) = 0$
- $\mathcal{D}(z)\vec{H}_F(z) \mathcal{D}'(z)\vec{H}_G(z) = 0$

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 solutions.

Hence, at an equilibrium point, either  $\mathcal{D}(z)$  or  $\mathcal{D}'(z)$  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  $\{\mathcal{D} = \mathcal{D}' = 0\}$ , we may assume that the point z is neither O nor N. Hence, F(z) and G(z) are non-zero, and so  $\mathcal{D}(z) = \mathcal{D}'(z) = 0$ .

**Eigenvalues of the linearization** We consider the eigenvalue decomposition of the matrix

$$\mathcal{A} = \operatorname{Jac}(\mathcal{D}\vec{H}_F - \mathcal{D}'\vec{H}_G)$$

on equilibrium point, given as the union of points satisfying Eq. (28) and (29).

**Solutions of Eq.** (28) If  $z_1 = z_2$ , the matrix  $\mathcal{A}$  has rank 2, and its characteristic polynomial is

$$T^{6}\left(T^{2} - \frac{\varepsilon(\varepsilon-1)(\varepsilon-2)\left(8(\Gamma-\gamma)^{2}P_{y_{1}}y_{1}y_{2} + P_{z_{2}}(2P_{z_{2}} - \gamma)P_{y_{2}}\right)}{2P_{y_{1}}(\Gamma-\gamma)y_{1}}T + \left(\frac{(\varepsilon-1)\gamma P_{y_{1}}(2\Gamma-\gamma)}{y_{1}}\right)^{2}\right)$$

**Solutions of Eq.** (29) If  $y_2 = -\frac{y_2 p_{z_1}}{(\varepsilon - 1)p_{z_2}}$ , the matrix  $\mathcal{A}$  has rank 2, and its characteristic polynomial is

$$T^{6}\left(T^{2}-4\varepsilon y_{2}(\varepsilon-1)(\varepsilon-2)(\Gamma-\gamma)T+\left(\frac{2\gamma(z_{1}-z_{2})(\varepsilon-1)^{2}(2\Gamma-\gamma)(\Gamma-\gamma)}{p_{z_{1}}}\right)^{2}\right).$$

The discriminant of the degree 2 factor factors as

$$\frac{16(\Gamma-\gamma)^2(\varepsilon-1)^2\left(a_6(z_1,y_2)-b_6(z_1,z_2)\right)\left(a_6(z_1,y_2)+b_6(z_1,z_2)\right)}{p_{z_1}^2}$$

with

$$a_6(z_1, y_2) = \varepsilon(\varepsilon - 2) P_{z_1} y_2$$
  
 $b_6(z_1, z_2) = (2\Gamma - \gamma)(\varepsilon - 1) \gamma(z_1 - z_2)$ 

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.

# 3.4 Algebraic computations for the contrast problem

# 3.4.1 Subcase of water (2 projective parameters)

We consider the set of singularities of the hypersurface  $\{D=0\}$  lying in the Bloch ball. We show in the following that these singularities can be computed by means of explicit rational expressions. These explicit solutions show how complex is the structure of the algebraic variety corresponding to the singularities of  $\{D=0\}$ . The number of real singularities can be computed directly without knowing the exact expressions by an algorithm ([9]) using in an optimized way the Cylindric Algebraic decomposition algorithm.

Statement of the main results We keep the same letters for the sake of simplification. The new coordinates of the center O of the Bloch ball are O:(0,-1,0,-1). We consider the set of polynomials  $S = \{D, D_{y_1}, D_{y_2}, D_{z_1}, D_{z_2}\}$ . We recall that we assume that  $2\Gamma_2 \ge \gamma_2 > 0$ , and that  $(\gamma_1, \Gamma_1) \ne (\gamma_2, \Gamma_2)$ . We normalize by means of  $\Gamma_1 = \gamma_1 = 1$  (water), leaving free the two parameters  $\Gamma_2$ ,  $\gamma_2$  corresponding to the second matter. Under this normalization, the hypothesis becomes  $(\gamma_2, \Gamma_2) \ne (1, 1)$ . Let  $\Pi$  be the plane  $\{y_1 = y_2 = 0\}$ .

The following theorems were proved in [9] using a real roots classification algorithm.

## **Theorem 3.** Consider the 9 polynomials:

$$\begin{split} f_1 &= \Gamma_2 - 1 \\ f_2 &= 3\Gamma_2 - 2\gamma_2 - 1 \\ f_3 &= 3\Gamma_2^2 - 5\Gamma_2\gamma_2 + \gamma_2^2 + 2\Gamma_2 - 2\gamma_2 + 1 \\ f_4 &= 2\Gamma_2^2 - 5\Gamma_2\gamma_2 + 2\gamma_2^2 - 2\Gamma_2 + 3\gamma_2 \\ f_5 &= 2\gamma_2^3 - (3\Gamma_2 + 11)\gamma_2^2 + \left(9\Gamma_2 + 6 - 3\Gamma_2^2\right)\gamma_2 + 2\Gamma_2\left(\Gamma_2 + 2\right)\left(\Gamma_2 - 1\right) \\ f_6 &= \Gamma_2 - 2\gamma_2 + 1 \\ f_7 &= 2\Gamma_2 - \gamma_2 - 1 \\ f_8 &= \gamma_2 - 2 + \Gamma_2 \\ f_9 &= 2\Gamma_2^2 - 5\Gamma_2\gamma_2 + 2\gamma_2^2 + 1. \end{split}$$

The zeroes of their product divide the subset of  $\mathbb{R}^2$  defined by  $2\Gamma_2 > \gamma_2 > 0$  into connected components where the cardinality of  $V_{\mathbb{R}} \cap \pi^{-1}(\gamma_2, \Gamma_2)$  is constant.

Using a Cylindrical Algebraic Decomposition, one can make this statement more precise by expliciting this cardinality. Let  $\psi : (y_1, z_1, y_2, z_2) \mapsto (-y_1, z_1, -y_2, z_2)$  be the symmetry fixing  $\Pi = \{y_1 = y_2 = 0\}$ , and let us consider the semi-algebraic sets (see Fig. 6):

$$\begin{split} \mathcal{G}_{1}^{-} &= \{0 < \gamma_{2} < 2\Gamma_{2}, \Gamma_{2} < 1, f_{2} > 0\}, \\ \mathcal{G}_{1}^{+} &= \{0 < \gamma_{2} < 2\Gamma_{2}, \Gamma_{2} > 1, f_{2} < 0, f_{4} > 0\}, \\ \mathcal{G}_{2}^{-} &= \{0 < \Gamma_{2} < 1, f_{6} > 0, f_{3} < 0\}, \\ \mathcal{G}_{2}^{+} &= \{\Gamma_{2} > 1, f_{6} < 0, f_{5} > 0\}, \\ \mathcal{G}_{1}^{+} &= \mathcal{G}_{1}^{+} \cup \mathcal{G}_{1}^{-}, \\ \mathcal{G}_{2} &= \mathcal{G}_{2}^{+} \cup \mathcal{G}_{2}^{-}, \\ \mathcal{G} &= \mathcal{G}_{1} \cup \mathcal{G}_{2}. \end{split}$$

**Theorem 4.** For all  $(\gamma_2, \Gamma_2)$  such that  $2\Gamma_2 > \gamma_2 > 0$ , the center O of the Bloch ball B is a singularity of  $\{D = 0\}$ . And if  $(\gamma_2, \Gamma_2) \in \mathcal{G}$ , there exist at most two other singularities in the interior of the Bloch ball:

- 1. if  $(\gamma_2, \Gamma_2) \in \mathcal{G}_1$  there is one other singularity lying on  $\Pi \cap \mathcal{B}$ ;
- 2. if  $(\gamma_2, \Gamma_2) \in \mathcal{G}_2$  there are two other singularities in  $\mathcal{B}$ ,  $\psi$ -symmetric, outside  $\Pi$ .

The configuration is illustrated in Figs. 6 and 7. Observe that the number of singularities inside  $\mathcal{B}$  is an invariant of the contrast problem. Two of the pairs of biological matters studied in [7], water-cerebrospinal fluid (normalized parameters  $[\gamma_2 = \frac{5}{4}, \Gamma_2 = \frac{25}{3}]$ ) and water-fat (normalized parameters  $[\gamma_2 = \frac{25}{2}, \Gamma_2 = 25]$ ) correspond to points outside  $\mathcal{G}$ , and their invariant is 1 in both cases (see Fig. 7).

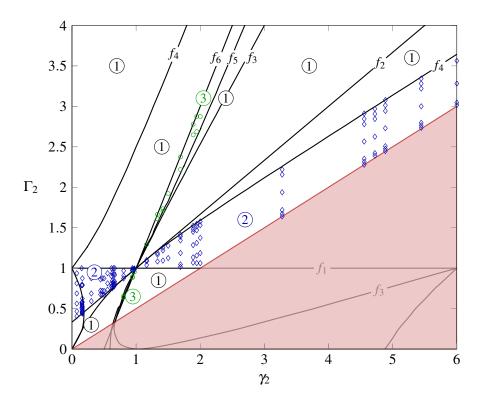


Figure 6: Curves involved in the definition of the semi-algebraic set  $\mathcal{G}$ . The blue (*resp.* green) sample points correspond to points in  $\mathcal{G}_1^- \cup \mathcal{G}_1^+$  (*resp.*  $\mathcal{G}_2^- \cup \mathcal{G}_2^+$ ). The circled numbers in each area correspond to the number of singularities in B for parameters in the area. Parameters in the red area are physically irrelevant.

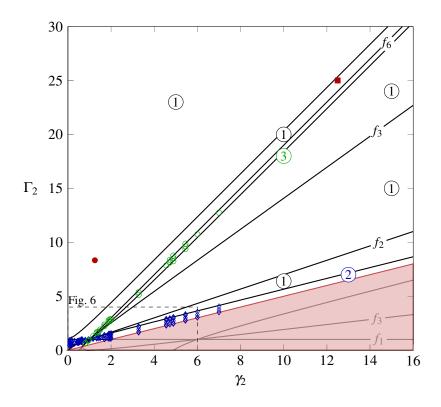


Figure 7: Positions of the parameters corresponding to the pairs water-cerebrospinal fluid (red circle) and water-fat (red square) and the set  $\mathcal{G}$  (with the same conventions as in Fig. 6). For both these pairs, there is only 1 singularity in  $\mathcal{B}$ .

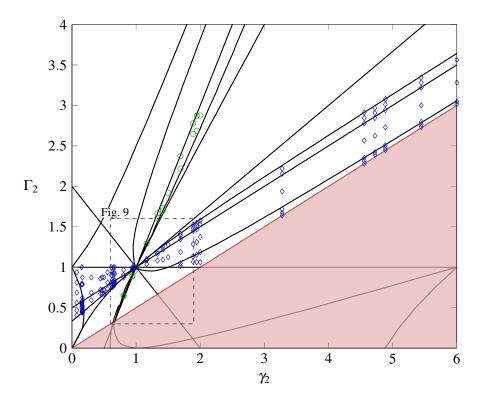


Figure 8: The curves involved in the decomposition of the parameter space (with the same conventions as in Fig. 6).

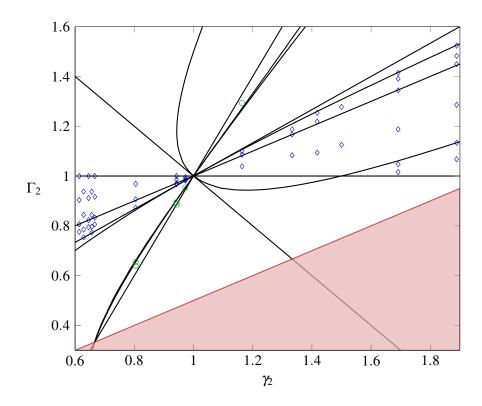


Figure 9: Magnification of Fig. 8 near (1,1).

We can give a more precise description of those points. Assuming that  $(\gamma_2 - \Gamma_2)(2\Gamma_2 - \gamma_2 - 2) \neq 0$ , we define a point (with the coordinates  $(y_1, z_1, y_2, z_2)$ ):

$$\Omega:\left(0,\frac{\Gamma_2-1}{\Gamma_2-\gamma_2},0,\frac{\left(1-\Gamma_2\right)\left(2\,\Gamma_2-\gamma_2\right)}{\left(2\,\Gamma_2-\gamma_2-2\right)\left(\Gamma_2-\gamma_2\right)}\right)$$

In the same coordinates system, O:(0,-1,0,-1) is the center of the Bloch ball, while N:(0,0,0,0) is the North pole.

Let

$$\delta(\gamma_2, \Gamma_2) = (\Gamma_2 - 1)(2\Gamma_2 - \gamma_2)(\Gamma_2 - 2\gamma_2 + 1)(\Gamma_2 - 2\gamma_2 + 2)$$
(30)

and

$$\eta(\gamma_2, \Gamma_2) = (5\Gamma_2 - 4\gamma_2 + 1)(\Gamma_2 - \gamma_2)(\Gamma_2 + \gamma_2 - 1)(\Gamma_2 - 2\gamma_2 + 2).$$

If  $\eta(\gamma_2, \Gamma_2) \neq 0$  we denote by  $\Omega_a^+$  (resp.  $\Omega_a^-$ ) the points of coordinates  $(y_1, z_1, y_2, z_2)$  (resp.  $(-y_1, z_1, -y_2, z_2)$ ), with:

$$\begin{split} y_1 &= \frac{\sqrt{3}\sqrt{\delta(\gamma_2, \Gamma_2)}}{(5\Gamma_2 - 4\gamma_2 + 1)(\Gamma_2 - \gamma_2)} \\ z_1 &= \frac{(\gamma_2 + \Gamma_2 - 1)(\Gamma_2 - 1)}{(5\Gamma_2 - 4\gamma_2 + 1)(\gamma_2 - \Gamma_2)} \\ y_2 &= \frac{\gamma_2\sqrt{3}\sqrt{\delta(\gamma_2, \Gamma_2)}}{2(\gamma_2 + \Gamma_2 - 1)(\Gamma_2 - 2\gamma_2 + 2)(\Gamma_2 - \gamma_2)} \\ z_2 &= \frac{(2\Gamma_2 - \gamma_2)(\Gamma_2 - 1)}{2(\gamma_2 + \Gamma_2 - 1)(\gamma_2 - \Gamma_2)}, \end{split}$$

**Theorem 5.** Generically on  $(\gamma_2, \Gamma_2)$ , the following points are singular points of  $\{D = 0\}$  in the subcase of water:

- O (always in the interior of the Bloch ball);
- $\Omega \ if (\gamma_2 \gamma_2)(2\Gamma_2 \gamma_2 2) \neq 0;$
- $\Omega_a^+$  and  $\Omega_a^-$  if  $\eta(\gamma_2, \Gamma_2) \neq 0$  and which are real if  $\delta(\gamma_2, \Gamma_2) > 0$ .

The remainder of this section will be devoted to the proof of Theorem 5. Beforehand, however, let us state this last theorem, which makes the connection between the results of Theorems 4 and 5, and in particular implies that generically, the points listed in Theorem 5 are the only singularities of  $\{D=0\}$  in the Bloch ball.

**Theorem 6.** Generically on  $(\gamma_2, \Gamma_2)$ , the singularities of  $\{D = 0\}$  inside the Bloch ball in the case of water are:

- O;
- $\Omega$  if and only if  $(\gamma_2, \Gamma_2) \in \mathcal{G}_1$ ;
- $\Omega_a^+$  and  $\Omega_a^-$  are real and inside the Bloch ball if and only if  $(\gamma_2, \Gamma_2) \in \mathcal{G}_2$ .

Remark 2. Note that generically the number of singularities of  $\{D=0\}$  in the Bloch ball is at most 3, because the domains  $\mathcal{G}_1$  and  $\mathcal{G}_2$  are disjoint.

*Proof.* The point  $\Omega$  is within the ball if and only if its coordinates satisfy

$$-2 < z_1, z_2 < 0.$$

The inequalities in  $z_1$  are equivalent to the disjonction of these two sets of inequalities

$$\begin{cases} \Gamma_2 - 1 > 0 \\ \Gamma_2 < \gamma_2 \\ 3\Gamma_2 - 2\gamma_2 - 1 = f_2 < 0 \end{cases} \quad \text{or} \quad \begin{cases} \Gamma_2 - 1 < 0 \\ \Gamma_2 > \gamma_2 \\ f_2 > 0 \end{cases}$$

Note that in each case, the first and third condition imply the second one.

In the first case, since  $\Gamma_2 > \gamma_2$ , the inequalities  $0 > z_2 > -2$  are equivalent to

$$\begin{cases} 2\Gamma_2 - \gamma_2 - 2 < 0 \\ f_4 < 0 \end{cases}$$

Both inequalities are satisfied by assumption, recovering the defining inequalities of  $\mathcal{G}_1^-$ . In the second case  $(\Gamma_2 > \gamma_2)$ , the inequalities  $0 > z_2 > -2$  are equivalent to

$$\begin{cases} 2\Gamma_2 - \gamma_2 - 2 > 0 \\ f_4 > 0 \end{cases}$$

The first inequality is satisfied whenever the second is, so have recoved the defining inequalities of  $\mathcal{G}_1^+$ , proving the first point.x

The second case is studied in the same way.

#### Proof of Theorem 5

*Proof.* We consider separately those solutions which are on the symmetry plane  $\{y_1 = y_2 = 0\}$  (Proposition 10) and those which are not (Proposition 11). The result is an immediate consequence of these lemmas.

Solutions on the symmetry plane  $\{y_1 = y_2 = 0\}$ 

**Proposition 10.** The singularities of D which are on the symmetry hyperplane  $\Pi$  are:

- if  $(2\Gamma_2 \gamma_2)(\Gamma_2 \gamma_2)(2\Gamma_2 \gamma_2 2) \neq 0$ : the center O of B, and the point  $\Omega$ ;
- if  $2\Gamma_2 = \gamma_2$ : O,  $\Omega$  and N:
- if  $\Gamma_2 \neq 1$ , and  $(\Gamma_2 \gamma_2)(2\Gamma_2 \gamma_2 2) = 0$ : O.
- if  $\Gamma_2 = 1$  and  $\gamma_2 \neq 1$ : O and  $\Omega = N$ .

*Proof.* We split the proof in seven cases, finding the solutions such that  $z_2 = -1$ ,  $\Gamma_2 = 1$ ,  $z_2 = 0$ ,  $2\Gamma_2 = \gamma_2$ ,  $\Gamma_2 = \gamma_2$ ,  $2\Gamma_2 - \gamma_2 - 2 = 0$ , and the remaining general case.

- 1. Case  $z_2 = -1$ . We compute a basis of  $S \cup \{y_1, y_2, z_2 + 1\}$  with the lexicographical order  $y_1 \gg y_2 \gg z_1 \gg z_2 \gg \gamma_2 \gg \Gamma_2$ . The polynomial  $\gamma_2^2(z_1 + 1)$  is present, and since  $\gamma_2 \neq 0$ , we get  $z_2 = -1$  and derive directly that the unique solution is O.
- 2. Case  $2\Gamma_2 = \gamma_2$ . We compute a basis of  $S \cup \{y_1, y_2, 2\Gamma_2 \gamma_2\}$  with the same order  $y_1 \gg y_2 \gg z_1 \gg z_2 \gg \gamma_2 \gg \Gamma_2$ , and the result follows immediately.
- 3. Case  $\Gamma_2 = 1$ . We found in the basis of  $S \cup \{y_1, y_2, \Gamma_2 1\}$  with order  $y_1 \gg y_2 \gg z_1 \gg z_2 \gg \gamma_2 \gg \Gamma_2$ , the polynomial  $z_2 \gamma_2^2 (\gamma_2 1)^2 (z_2 + 1)$ . The case  $z_2 + 1$  is already treated, while the case  $z_2 = 0$  is excluded. The case  $z_2 = 1$  is also excluded, since that would mean  $(\gamma_2, \Gamma_2) = (1, 1)$ . In the case  $z_2 = 0$ , the polynomial  $2\gamma_2^3 z_1 (\gamma_2 2)$  appears in the basis of  $S \cup \{y_1, y_2, \Gamma_2 1, z_2\}$ . Provided that  $z_2 \neq 2$ , this implies  $z_1 = 0$  hence the north pole  $z_1 = 0$  is the unique solution. Note that for  $z_2 = 1$ ,  $z_2 = 1$ ,
- 4. Case  $z_2=0$  and  $\Gamma_2 \neq 1$ . We compute a basis with the lexicographic  $y_1 \gg y_2 \gg z_1 \gg z_2 \gg \gamma_2 \gg \Gamma_2$ . The first polynomial is  $(\Gamma_2 1)^2 (2\Gamma_2 \gamma_2)$ . Since  $2\Gamma_2 \gamma_2 = 0$  and  $\Gamma_2 = 0$  are already known, there is no extra solution.
- 5. Case  $\Gamma_2 = \gamma_2$  and  $\Gamma_2 \neq 1$ . The basis contains the polynomial  $\Gamma_2(\Gamma_2 1)^2(z_2 + 1)$ , which leads to  $z_2 = -1$  (already discussed).
- 6. Case  $2\Gamma_2 \gamma_2 2 = 0$  and  $\Gamma_2 \neq 1$ . The basis contains the polynomial  $(\Gamma_2 1)^2(z_2 + 1)$ , as in the previous case.
- 7. Let us finally search the solutions such that  $z_2(z_2+1)(\gamma_2-\Gamma_2)(2\Gamma_2-\gamma_2-2)\neq 0$ . We consider the system

$$S \cup \{y_1, y_2, z_2\xi_1 - 1, (z_2 + 1)\xi_2 - 1, (\Gamma_2 - \gamma_2)\lambda_1 - 1, (2\Gamma_2 - \gamma_2 - 2)\lambda_2 - 1\},\$$

and compute a basis with order  $y_1 \gg y_2 \gg z_1 \gg z_2 \gg \xi_1 \gg \xi_2 \gg \gamma_2 \gg \Gamma_2 \gg \lambda_1 \gg \lambda_2$ . The last four polynomials of the basis are:

$$\Gamma_2 \lambda_1 + 2 \lambda_2 \lambda_1 + \lambda_1 - 2 \lambda_2 + z_2, -\Gamma_2 \lambda_1 + \lambda_1 + z_1, y_2, y_1.$$

This implies that there is at most one solution under the above conditions. By plugging into the system, one verifies that  $\Omega$  is a solution satisfying these conditions, hence it is the unique solution.

Solutions outside the symmetry plane  $\{y_1 = y_2 = 0\}$ 

**Proposition 11.** The solutions of S outside  $\Pi$  are given by means of

- 1. a two parameters family of  $\Pi$ -symmetric points  $(\Omega_a^+(\gamma_2, \Gamma_2), \Omega_a^-(\gamma_2, \Gamma_2))$  which can be either real or complex according to the sign of  $\delta(\gamma_2, \Gamma_2)$ ;
- 2. non-generically on the parameters, two one-parameter families of complex points.

These one-parameter families are explicitly given by:

1. if  $\gamma_2 = 2 - \Gamma_2$ ,  $\Gamma_2 \neq 1$ ,  $\Gamma_2 \neq 2$ ,  $z_1 = \frac{1}{4} \frac{\Gamma_2}{1 - \Gamma_2}, z_2 = \frac{1}{4} \frac{3\Gamma_2 - 2}{1 - \Gamma_2}, t = \frac{3\Gamma_2 - 2}{\Gamma_2 - 2}, Y_2 = -\frac{1}{16} \frac{(\Gamma_2 - 2)^2}{(\Gamma_2 - 1)^2}$ 

2. if  $2\Gamma_2^2 - 5\gamma_2\Gamma_2 + 2\gamma_2^2 + 1 = 0$ , it means that  $(\gamma_2, \Gamma_2)$  belong to a hyperbola which can be parameterized as

$$\Gamma_2 = \frac{\sqrt{2} (5 u^2 + 6u + 5)}{6 (1 - u^2)}$$

$$\gamma_2 = \frac{2 \sqrt{2} (u^2 + 1)}{3 (1 - u^2)}$$

If  $|u| \neq 1$  and  $|u+3| \neq 2\sqrt{2}$ , the corresponding solutions in  $(Y_2,t,z_1,z_2)$  are given by

$$Y_{2} = -\frac{8(3+2\sqrt{2})(u+1)^{2}(u^{2}+1)^{2}}{(2\sqrt{2}-u-3)^{2}(u+3+2\sqrt{2})^{4}}$$

$$t = \frac{3(u-1)^{2}}{4(u^{2}+1)}$$

$$z_{1} = \frac{(3\sqrt{2}+8)(-23u^{3}+(14\sqrt{2}-45)u^{2}+(-36\sqrt{2}-65)u-10\sqrt{2}-27)}{46(u+3-2\sqrt{2})(u+3+2\sqrt{2})^{2}}$$

$$z_{2} = \frac{(2\sqrt{2}+1)(3-u+2\sqrt{2})(4\sqrt{2}-7u-9)(u+1)}{7(u+3-2\sqrt{2})(u+3+2\sqrt{2})^{2}}.$$

Observe that since  $2\Gamma_2 > \gamma_2$  the first solution has to be considered provided that  $0 < \gamma_2 < \frac{4}{3}$  and  $\Gamma_2 > \frac{2}{3}$ .

# Proof of Proposition 11

*Proof.* The case  $y_2 = 0, y_1 \neq 0$  is treated in Lemma 11. We assume now that  $y_2 \neq 0$ .

We transform the system  $S = \{D, D_{y_1}, D_{y_2}, D_{z_1}, D_{z_2}\}$  by means of the introduction of the new variables (t,h) such that  $y_1 = ty_2$ , and  $\gamma_2 = h\Gamma_2$ . After simplification by  $y_2$ , one observes that the powers of  $y_2$  in the simplified system all all even. So we perform the change of variables  $Y_2 = y_2^2$ , and denote by  $\tilde{S}$  this new set of polynomials in the variables  $(h, \Gamma_2, z_1, z_2, t, Y_2)$ .

We first compute a Gröbner basis, considering the polynomials in  $\mathbb{Q}(\gamma_2, \Gamma_2)[y_1, z_1, y_2, z_2]$  with a lexicographical order, and find directly the rational expressions

$$\begin{split} z_1 &= \frac{\left(1 - \left(1 + h\right)\Gamma_2\right)\left(\Gamma_2 - 1\right)}{\left(4\Gamma_2 h - 5\Gamma_2 - 1\right)\Gamma_2\left(h - 1\right)} \\ z_2 &= \frac{1}{2} \frac{\left(\Gamma_2 - 1\right)\left(h - 2\right)}{\left(1 - \left(1 + h\right)\Gamma_2\right)\left(h - 1\right)} \\ Y_2 &= \frac{3}{4} \frac{h^2\Gamma_2\left(\Gamma_2 - 1\right)\left(h - 2\right)\left(2\Gamma_2 h - \Gamma_2 - 1\right)}{\left(2\Gamma_2 h - \Gamma_2 - 2\right)\left(\Gamma_2 h + \Gamma_2 - 1\right)^2\left(h - 1\right)^2} \\ t &= 2 \frac{\left(\Gamma_2 h + \Gamma_2 - 1\right)\left(2\Gamma_2 h - \Gamma_2 - 2\right)}{\left(4\Gamma_2 h - 5\Gamma_2 - 1\right)h\Gamma_2} \end{split}$$
 (SoluGen)

which correspond to the points  $\Omega_a^+, \Omega_a^-$ . These points are symmetric with respect to  $\Pi$  and real if and only if

$$(\Gamma_2 - 1)(2\Gamma_2 - h\Gamma_2)(\Gamma_2 - 2h\Gamma_2 + 1)(\Gamma_2 - 2h\Gamma_2 + 2) \ge 0$$

Recall that  $h\Gamma_2 = \gamma_2$ , so we have recovered the definition of  $\delta$  (30). We define

$$\bar{\eta}(\Gamma_2, h) = (4\Gamma_2 h - 5\Gamma_2 - 1)(h - 1)(2\Gamma_2 h - \Gamma_2 - 2)(1 - (1 + h)\Gamma_2).$$

the lowest common multiple of the denominators of the solutions, so that those paired solutions are valid on a Zariski open set (including at least the complementary of  $\{(\gamma_2, \Gamma_2) \text{ s.t. } \bar{\eta}(\Gamma_2, h) = 0\}$ . Let us search extra solutions.

We compute a basis of  $\tilde{S}$  with respect to the elimination order  $Y_2 > z_2 > z_1 \gg t > \Gamma_2 > h$ . In this basis, one gets the polynomial:

$$h^{2}(h-2)^{2}(\Gamma_{2}-1)\theta_{1}(h,t)\theta_{2}(h,t,\Gamma_{2})\theta_{3}(h,t,\Gamma_{2})$$

with

$$\begin{split} \theta_{1}(h,t) = & 1 + (t-2)h \\ \theta_{2}(h,t,\Gamma_{2}) = & \Gamma_{2}^{2}(t-1)h^{2} - \Gamma_{2}\left(5\Gamma_{2}t + 2\Gamma_{2} + t - 8\right)h + 2\left(\Gamma_{2} + 2\right)\left(\Gamma_{2} - 1\right) \\ \theta_{3}(h,t,\Gamma_{2}) = & -2h(2h-3)t^{2} \\ & + \left(\left(4\Gamma_{2} + 4\right)h^{2} + \left(-6\Gamma_{2}^{2} + 10\Gamma_{2} - 16\right)h - 3\left(\Gamma_{2} + 2\right)\left(\Gamma_{2} - 1\right)\right)t \\ & + 4\left(2\Gamma_{2} - 1\right)\left(\Gamma_{2} - 2\right)h^{2} + \left(-2\Gamma_{2}^{2} + 6\Gamma_{2} + 2\right)h - \left(\Gamma_{2} - 1\right)\left(\Gamma_{2} - 10\right)\Gamma_{2} \end{split}$$

The case h = 0 is irrelevant. The case h = 2 (resp.  $\Gamma_2 = 1$ ) is treated in Lemma 9 (resp. Lemma 7). We split the remainder of the study in three parts.

1. We assume that  $Y_2 \neq 0$  and introduce its inverse  $\xi_1$ . We compute a basis of the system  $\tilde{S} \cup \{\theta_1(h,t), Y_2 \xi_1 - 1\}$  with respect to the DRL order  $\xi_1 > Y_2 > z_2 > z_1 > t \gg \Gamma_2 > h$ . The first polynomial is:

$$(h-1)(2h\Gamma_2-7\Gamma_2+4)(2\Gamma_2^2h^2-5\Gamma_2^2h+2\Gamma_2^2+1)$$

while in the second polynomial, the factor  $\Gamma_2 - 1$  appears. The case h = 1 (resp.  $\Gamma_2 = 1$ ) is treated in Lemma 8 (resp. Lemma 7).

We assume that  $\Gamma_2 - 1 \neq 0$  and introduce its inverse  $\lambda_1$ . We compute a basis of

$$\tilde{S} \cup \{\theta_1(h,t), Y_2\xi_1 - 1, 2h\Gamma_2 - 7\Gamma_2 + 4, (\Gamma_2 - 1)\lambda_1 - 1\}$$

with respect to the lexicographical order  $\lambda_1 \gg \xi_1 \gg z_2 \gg z_1 \gg t \gg Y_2 \gg h \gg \Gamma_2$ , and this leads to a unique solution which is exactly the general solution evaluated at  $\gamma_2 = \frac{7}{2}\Gamma_2 - 2$ .

We consider the polynomial  $2\Gamma_2^2h^2 - 5h\Gamma_2^2 + 2\Gamma_2^2 + 1$  corresponding to an hyperbola, admitting the parameterization given in the statement. We compute a basis of

$$\tilde{S} \cup \{\theta_1(h,t), Y_2 \xi_1 - 1, 2\Gamma_2^2 h^2 - 5h\Gamma_2^2 + 2\Gamma_2^2 + 1, (\Gamma_2 - 1)\lambda_1 - 1\}$$

with respect to the lexicographical order  $\lambda_1 \gg \xi_1 \gg z_2 \gg z_1 \gg t \gg Y_2 \gg h \gg \Gamma_2$ . This leads directly to the solution item 2 of the statement of Prop. 11.

2. We consider the system  $\tilde{S} \cup \{\theta_3(h,t,\Gamma_2)\}$ . The first polynomial of a basis with respect to the order  $\xi_1 \gg z_2 \gg z_1 \gg t \gg Y_2 \gg \Gamma_2 \gg h$  is:

$$(h-1)(h-2)(2h-3)\theta_4(Y_2,h,\Gamma_2)$$

with

$$\theta_4(Y_2, h, \Gamma_2) = 4 (h-1)^2 (2 h \Gamma_2 - \Gamma_2 - 2) (h \Gamma_2 + \Gamma_2 - 1)^2 Y_2$$
$$-3 \Gamma_2 h^2 (h-2) (\Gamma_2 - 1) (2 h \Gamma_2 - \Gamma_2 - 1)$$

The cases h = 1, h = 2,  $h = \frac{3}{2}$  are studied in Lemmas 8, 9, 10.

If  $\bar{\eta}(\Gamma_2, h) \neq 0$ , the system is linear in  $(z_1, z_2, h, Y_2, \xi_1)$  and non-singular, and there is a unique solution, corresponding to (SoluGen).

To examine the case where  $\bar{\eta}(\Gamma_2, h) = 0$ , we compute a Gröbner basis of

$$\tilde{S} + \langle \theta_3, \theta_4, \xi_1 Y_2 - 1, u(h-1)(h-2)(2h-3) - 1 \rangle$$

for the elimination order  $u > \xi_1 \gg z_2 > z_1 > t > Y_2 > G_2 > h$ , and we find that this basis is  $\{1\}$ , so there is no extra solution.

3. We consider the system

$$\tilde{S} \cup \{\theta_2(h,t,\Gamma_2), \theta_1(h,t,\Gamma_2)\mu_1 - 1, \theta_3(h,t,\Gamma_2)\mu_3 - 1, Y_2\xi_1 - 1\}$$

and compute a basis with respect to the elimination order  $\mu_1 > \mu_3 > \xi_1 > Y_2 > z_1 > z_2 > t \gg \Gamma_2 > h$ .

The first polynomial of the basis is  $\Gamma_2 h + \Gamma_2 - 2$ , and the corresponding solutions are given in Lemma 6.

We should mention that the choice of a suitable ordering is a key feature, in order to achieve the computation of a basis, and to find a factorized polynomial in it. When a factorization is found, the difficulty of the computations with each factor falls dramatically. In the last case of the proof, it is absolutely necessary to saturate in order to isolate the polynomial  $\Gamma_2 h + \Gamma_2 - 2$ . On the other hand, dealing with too many saturations increases the number of variables and can lead to a practical impossibility of computing a basis.

**Lemma 6.** Provided that  $h\Gamma_2 + \Gamma_2 - 2 = 0$ , the system  $\tilde{S}$  has no other solution than those listed in Prop. 11.

*Proof.* The first polynomial of the basis of  $\tilde{S} \cup \{h\Gamma_2 + \Gamma_2 - 2\}$  with respect to the lexicographical order  $\Gamma_2 \gg z_1 \gg z_2 \gg t \gg Y_2 \gg h$  is  $(h-1)^2 (h-2)^2 F_1 F_2$  with  $F_1 = 4 (h+1)^2 Y_2 + 9 h^2$  and  $F_2 = 4 (h-1)^2 Y_2 + h^2$ . The case h=1 (resp. h=2) is treated in Lemma 8 (resp. Lemma 9).

1. If  $F_1 = 0$ , we compute a basis of

$$\tilde{S} \cup \{F_1, F_2F_I - 1, (h-1)h_I - 1, (h-2)h_{2I} - 1\}$$

with order  $h_I \gg h_{2I} \gg F_I \gg z_1 \gg z_2 \gg t \gg Y_2 \gg \Gamma_2 \gg h$ . The case  $\Gamma_2 = 2$  corresponding to  $\gamma_2 = 0$  is irrelevant. If  $9\Gamma_2 = 7$  there is no solution. The system is linear with respect to the variables  $z_1, z_2, t, Y_2$ , and has a unique solution provided that  $\Gamma_2 \neq 2$  and  $9\Gamma_2 \neq 7$ . This solution corresponds to (SoluGen) evaluated at  $\gamma_2 = 2 - \Gamma_2$ .

2. If  $F_2 = 0$ , we compute a basis of

$$\tilde{S} \cup \{F_2, F_1F_I - 1, (h-1)h_I - 1, (h-2)h_{2I} - 1\}$$

with order  $h_I \gg h_{2I} \gg F_I \gg z_1 \gg z_2 \gg t \gg Y_2 \gg \Gamma_2 \gg h$ . The case  $\Gamma_2 = 2$  means  $\gamma - 2 = 0$ , which is irrelevant. In the case  $\Gamma_2 = 1$  this last system has no solution, and is linear with respect to  $(z_1, z_2, t, Y_2)$ . If  $\Gamma_2 \neq 1$  and  $\Gamma_2 \neq 2$ , it has a unique solution given at item 1 of Lemma 11.

3. If  $F_1 = F_2 = 0$ , we compute a basis of

$$\tilde{S} + \langle h\Gamma_2 + \Gamma_2 - 2, F_1, F_2, u(h-1)(h-2) - 1 \rangle$$

for the order  $u \gg z_2 \gg z_1 \gg t \gg Y_2 \gg \Gamma_2 \gg h$ , and this basis gives a solution such that  $Y_2 = -\frac{1}{4}$ , so there is no real solution in  $y_2$ .

**Lemma 7.** If  $\Gamma_2 = 1$ , the solutions of S are

- 1. O
- 2. if  $\gamma_2 = 1$ , the plane  $\{y_2 = y_1, z_2 = z_1\}$ ,

3. if 
$$\gamma_2 = \frac{3}{2}$$
,

$$\left\{y_1 = \frac{y_2}{y_2^2 + 1}, z_1 = \frac{y_2^2}{y_2^2 + 1}, z_2 = 0, y_2 \in \mathbb{R}\right\}$$

*Proof.* A basis of  $S \cup \{\Gamma_2 - 1\}$  with respect to  $z_1 \gg z_2 \gg y_1 \gg y_2 \gg \gamma_2 \gg \Gamma_2$  gives the second polynomial

$$y_2^5 \gamma_2 (\gamma_2 - 1)^2 (2 \gamma_2 - 3)^2$$

and the solutions follow directly.

**Lemma 8.** Assume that  $\Gamma_2 = \gamma_2$ . If  $\gamma_2 = \Gamma_2 \neq 1$ , O is the unique solution of the system S. If  $\gamma_2 = \Gamma_2 = 1$  the solutions are the plane  $\{y_1 = y_2, z_1 = z_2, (y_2, z_2) \in \mathbb{R}^2\}$ .

*Proof.* The proof is straightforward from the computation of a basis of  $S \cup \{\Gamma_2 - \gamma_2\}$  with respect to the order  $y_1 \gg y_2 \gg z_1 \gg z_2 \gg \gamma_2 \gg \Gamma_2$ .

**Lemma 9.** If  $\gamma_2 = 2\Gamma_2$ ,  $\gamma_2 > 0$ , the solutions of S are:

- 1. the point O,
- 2. the circle

$$\left\{ \left( z_1 + 1 - \frac{1}{\gamma_2} \right)^2 + y_1^2 - \frac{1}{\gamma_2^2} = 0, \quad y_2 = z_2 = 0 \right\},\,$$

- 3. if  $\gamma_2 = \frac{4}{3}$ , the line  $\{(0, \frac{1}{2}, y_2, 0), y_2 \in \mathbb{R}\}$ ,
- 4. if  $\gamma_2 = \frac{2}{3}$ , the hyperbolic curve

$$\left\{ y_2 = -\frac{2y_1}{3z_1}, \quad z_2 = \frac{2-z_1}{3z_1}, \quad 2\left(z_1 + \frac{1}{2}\right)^2 - y_1^2 - \frac{1}{2} = 0, \quad z_1 \in \mathbb{R}^* \right\}.$$

The number of solutions is infinite and the dimension of the algebraic variety is 1.

*Proof.* We compute a basis for the system  $S \cup \{\gamma_2 - 2\Gamma_2\}$  with respect to the order  $y_1 > y_2 > z_1 > z_2 \gg \Gamma_2 > \gamma_2$ . The second polynomial of the basis is

$$z_2 \gamma_2 (3 \gamma_2 - 2)^2 (z_1 \gamma_2 - 2 z_2 + \gamma_2 - 2)$$

and a straightforward resolution gives the solutions listed in the statement.

**Lemma 10.** If  $h = \frac{3}{2}$  and  $\theta_3 = 0$   $\tilde{S}$  has the solutions:

- the solutions (SoluGen) for  $\gamma_2 = \frac{3}{2} \Gamma_2$ ;
- provided that  $\Gamma_2 = 1$ , the solutions listed in Lemma 7 item 3.

*Proof.* The second polynomial of a basis of  $\tilde{S} \cup \{\theta_3, 2h-3\}$  with respect to the order  $z_2 \gg z_1 \gg t \gg Y_2 \gg \Gamma_2 \gg h$  is

$$(\Gamma_2 - 1) \left( 4 (5\Gamma_2 - 2)^2 Y_2 + 27\Gamma_2 (2\Gamma_2 - 1) \right)$$

The case  $\Gamma_2 = 1$  is treated in Lemma 7, item 3. The basis of  $\tilde{S} \cup \{\theta_3, 2h - 3, 5\Gamma_2 - 2\}$  is  $\{1\}$ , hence the system has no solution. So we can express  $Y_2$  in terms of  $\Gamma_2$  and derive a unique solution which is (SoluGen) computed at  $\gamma_2 = \frac{3}{2}\Gamma_2$ .

**Lemma 11.** The system  $S \cup \{y_2\}$ , S has the following solutions:

- the solutions on  $\Pi$  given in Proposition 10;
- provided that  $\gamma_2 = 2\Gamma_2$ , the solutions given in Lemma 9 evaluated at  $y_1 = 0$ .

*Proof.* The second polynomial of a basis of  $S \cup \{y_2\}$  with respect to the order  $z_1 \gg z_2 \gg y_1 \gg y_2 \gg \gamma_2 \gg \Gamma_2$  is  $\gamma_2^2 y_1 (2\Gamma_2 - \gamma_2)$ . The case  $y_1 = 0$  (resp.  $\gamma_2 = 2\Gamma_2$ ) is treated in Proposition 10 (resp. Lemma 9).

#### 3.4.2 General case (3 projective parameters)

We conclude by giving classification results for the classification of the real singularities of  $\{D=0\}$  in the Bloch ball, in the general case. We now have 4 parameters  $(\gamma_1, \Gamma_1, \gamma_2, \Gamma_2)$ , which we may reduce to 3, normalizing by  $\gamma_1 = 1$ . The constraints are  $2\Gamma_1 \ge 1$ ,  $2\Gamma_2 \ge \gamma_2 > 0$ , and we exclude the case of a single spin  $(\Gamma_2, \gamma_2) = (\Gamma_1, 1)$ .

In this case, real roots classification algorithms [9] allow to compute a generic classification of the parameter space according to the number of singularities.

**Theorem 7** ([9]). Splitting the subset of  $\mathbb{R}^3$  defined by  $2\Gamma_2 \ge \gamma_2 > 0$  and  $2\Gamma_1 > 1$  into open subsets where the number of real singularities of  $\{D = 0\}$  in the Bloch ball in the fibers is constant, can be done by cutting out 12 irreducible surfaces, consisting of 5 planes, 3 quadrics, 2 surfaces of degree 9, and one of degree 14.

Computing a Cylindrical Algebraic Decomposition of the hypersurface defined by the product of these polynomials then allows to count the singularities. However, unlike the case of water, it is not possible to characterize the cells using only the signs of these polynomials.

**Theorem 8.** Generically on  $(\gamma_1, \Gamma_1, \gamma_2, \Gamma_2)$ , the number of singularities of  $\{D = 0\}$  in the Bloch ball can be 1, 2, 3, 4 or 5.

As an example, we show slices of this classification for  $\Gamma_1 = 531/26 \simeq 20.4$ , which corresponds to the *in vivo* experimental setting. All polynomials involved are represented on Figure 10. In this case, there can be 1, 2, 3 or 4 singularities, and the corresponding sample points are represented respectively with gray dots, blue diamonds, green circles and orange boxes.

In the experimental setting, the parameters for the second matter were  $\gamma_2 \simeq 0.89$  and  $\Gamma_2 \simeq 36.6$ . Scaling the graphic down to fit this point (Figure 11) shows that some of the polynomials can be eliminated for values in this order of magnitude. Several of the classification polynomials vanish at  $(\gamma_2, \Gamma_2) = (\gamma_1, \Gamma_1)$ , which is a natural singularity. One can also note that unlike in the case of water, the areas where there is only 1 singularity are quite far from the actual values of the parameters.

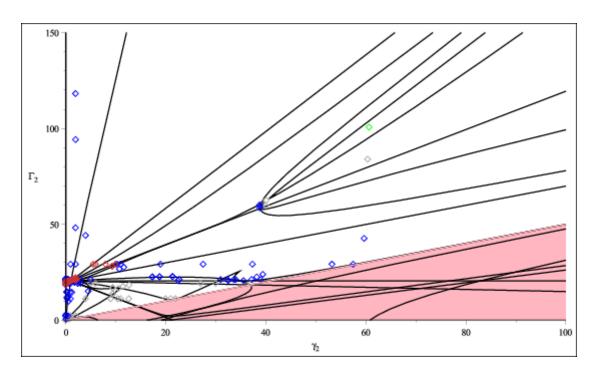


Figure 10: All polynomials involved in the classification in the in vivo case

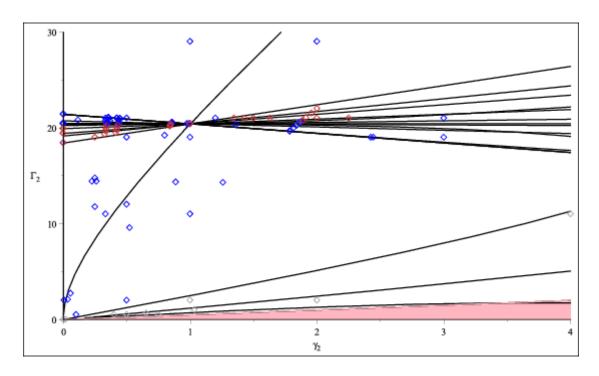


Figure 11: Classification regarding the second parameter in the *in vivo* case

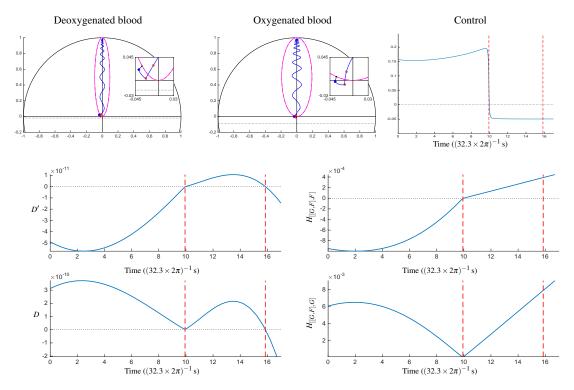


Figure 12: Singular flow for a pair of spins (deoxygenated blood and oxygenated blood), starting from a point near the colinearity locus of *F* and *G* 

## 4 Numerical simulations and discussion

## 4.1 Singular flows

Some numerical simulations of singular flows in the exceptional case. More precisely the role of  $D \cap D' = \{0\}$  is presented, showing that the studied invariants can have an impact on the trajectories. In each of Figures 12, 13 and 14, the vertical dashed red lines correspond to times at which D and D' vanish. They are illustrated by red dots on the trajectories.

Figure 12 corresponds to a contrast problem: spin 1 is deoxygenated blood, spin 2 is oxygenated blood, and we start the singular trajectory near the colinearity locus of F and G (the ellipse). In this case, when D and D' vanish for the first time, the trajectory has a singular point.

Figures 13 and 14 represent to a multisaturation problem with  $B_1$ -inhomogeneity, with spins of deoxygenated blood. The figures correspond to different starting points near the colinearity locus of F and [F,G] (the horizontal plane at altitude  $z_S$ ). On Figure 13, D and D' vanish only once, and the trajectory remains smooth. On Figure 14, D and D' vanish twice, the trajectory remains smooth. The first vanishing point is on the colinearity locus. The difference between the non smooth and smooth case is related to the simultaneous vanishing and non vanishing of  $H_{[[G,F],F]}$  and  $H_{[[G,F],G]}$ .

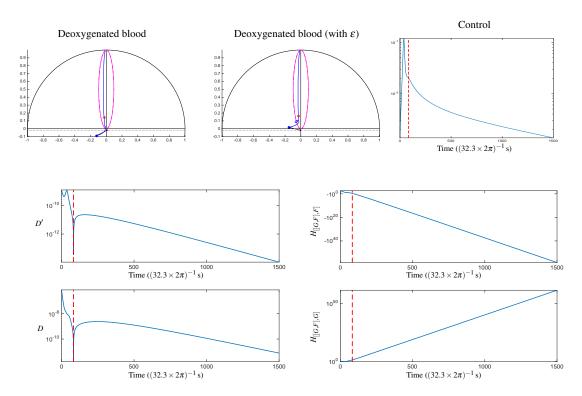


Figure 13: Singular flow for a pair of spins (deoxygenated blood with  $B_1$ -inhomogeneity), starting from a point near the colinearity locus of F and [F,G] (below the horizontal line for the first spin and above for the second)

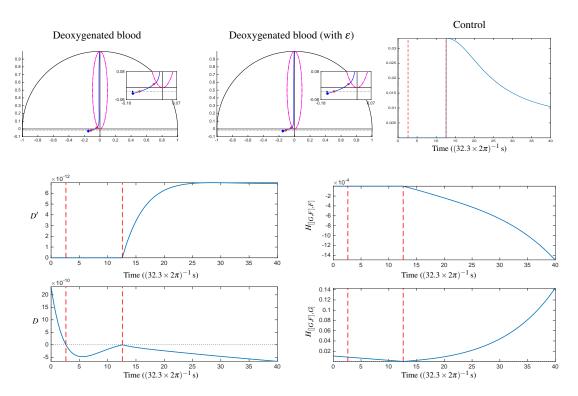


Figure 14: Singular flow for a pair of spins (deoxygenated blood with  $B_1$ -inhomogeneity), starting from a point near the colinearity locus of F and [F,G] (below the horizontal line for both spins)

## 4.2 Saturation problem

#### 4.2.1 Direct approach

We consider the optimal control problem where we saturate both spins in minimum time taking into account  $B_1$ -inhomogeneities. This problem can be formulated as

$$T_{min}^* = \min_{u(.)} t_f, \qquad |u| \le 1$$

$$\frac{dq}{dt} = V(q_1, q_2)$$

$$(q_1(0), q_2(0)) = q_i, \quad (q_1(t_f), q_2(t_f)) = q_f$$
(32)

where  $\varepsilon$  is the small rescaling parameter related to the  $B_1$ -inhomogeneity,  $q_i, q_f$  are respectively the initial and final state and  $V(q_1, q_2)$  is the vector field:

$$V(q_1, q_2) = (F(q_1) + uG(q_1)) \frac{\partial}{\partial q_1} + (F(q_2) + (1 - \varepsilon)uG(q_2)) \frac{\partial}{\partial q_2}$$

with F, G given by (5). In the sequel, numerical results are presented for the saturation of a single spin and of two spins, we let the reader adapt the formulations for the single spin case.

We present local solutions obtained by the Bocop software for the multisaturation problem in the Desoxygenated blood case  $(C_1)$ , the Oxygenated blood case  $(C_2)$ , the Cerebrospinal fluid case  $(C_3)$  and a Water "like" case  $(C_4)$ . The value of the parameters are:  $\varepsilon = 0.1$ ,  $q_i = (0,1,0,1)$ ,  $q_f = (0,0,0,0)$ . The time evolution of the state variables  $q = (q_1,q_2)$  and of the control variable u are represented in Fig.15, the optimal time is given in Table 1 and is compared with the optimal time of a saturation of a single spin.

Case	$T_1$	$T_2$	$\Gamma = (32.3 \times 2\pi T_1)^{-1}$	$\gamma = (32.3 \times 2\pi T_2)^{-1}$	$t_f$ (2 spins)	$t_f$ (1 spin)
$C_1$	1.35	0.05	$9.855 \times 10^{-2}$	$3.65 \times 10^{-3}$	44.769	42.685
$C_2$	1.35	0.2	$2.464 \times 10^{-2}$	$3.65 \times 10^{-3}$	113.86	110.44
$C_3$	2	0.3	$1.642 \times 10^{-2}$	$2.464 \times 10^{-3}$	168.32	164.46
$C_4$	0.05	0.05	$9.855 \times 10^{-2}$	$9.855 \times 10^{-2}$	8.7445	15.0237

Table 1: Cases treated numerically corresponding respectively to the Desoxygenated case, the Oxygenated case, the Cerebrospinal fluid case and a Water "like" case. The 7th (resp. 6th) column gives the local optimal time found by Bocop for the saturation of one spin (resp. two spins with  $B_1$ -inhomogeneity).

#### 4.2.2 Validation with the Linear Matrix Inequalities (LMI) approach

A crucial step is to check whether the local optimal times presented in Table 1 and obtained by Bocop for the saturation problem are globally optimal using moment/lmi techniques. These techniques provide lower bounds on the global optimal time  $T_{min}^*$  for the saturation problem of one spin or two spins (32). Note that in [6], the authors have investigated the contrast problem using these relaxations techniques.

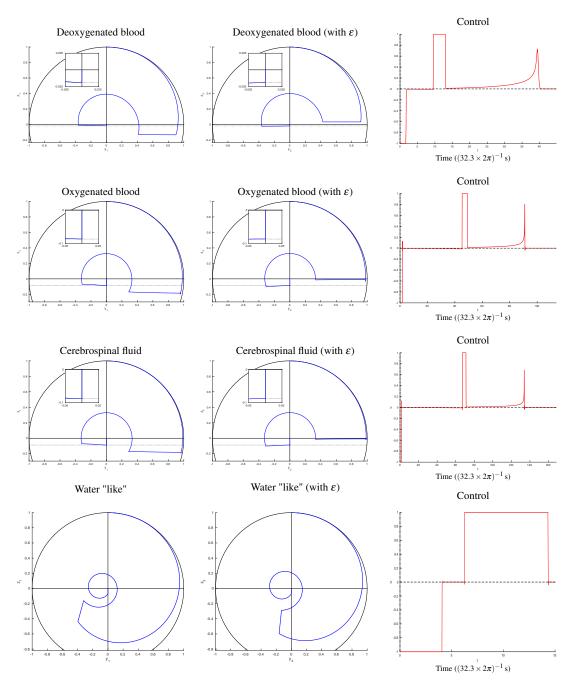


Figure 15: BC-extremal for the multisaturation problem with  $B_1$ -inhomogeneity in the cases  $C_1, C_2, C_3$  and  $C_4$ .

The moment approach is a global optimization method which relax 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* [31], this leads to an infinite dimensional LMI problem which can be truncated to a finite set of moments. The optimal value sequence associated with these truncated problems converges to the optimal value  $T_{min}^*$  of (32). Also, there is no efficient algorithm to solve the LMI hierarchy. We present an alternative formulation based on [12] which exploits the structure of the problem.

We first set up some 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 (32) can be embed into the linear program on measures:

$$T_{LP} = \inf_{\mu,\mu_f} \int d\mu_f$$

$$\int \left(\frac{\partial v}{\partial t}(t,q) + \frac{\partial v}{\partial q}(t,q)V(q)\right)d\mu$$

$$= \int v(\cdot,q_f)d\mu_f - v(q_i), \ \forall v \in \mathbb{R}[t,q],$$

$$\mu \in \mathcal{M}^+([0,T] \times Q \times U), \ \mu_f \in \mathcal{M}^+(Q_f)$$
(33)

where T is fixed,  $Q_f = [0, T]$ ,  $Q = B_2 \times B_2$  are admissible state sets and  $U = B_1$  is the admissible control set.

Given any admissible pair (q(.), u(.)) for (32), it corresponds a measure  $\mu$  admissible for (33) achieving the same cost, hence  $T_{min}^* \ge T_{LP}$ . Moreover, according to [25, Theorem 3.6 (ii)], there is no optimality gap and

$$T_{min}^* = T_{LP}$$
.

*Remark* 3. Since the dynamic and the Lagrangian are autonomous, the time variable can be removed and the LP problem becomes

$$T_{LP} = \inf_{\mu} \int d\mu$$
 
$$\int \frac{\partial v}{\partial q}(q)V(q)d\mu = v(q_f) - v(q_i), \quad \forall v \in \mathbb{R}[q],$$
  $\mu \in \mathcal{M}^+(O \times U).$ 

2nd formulation: Moment approach using modal occupation measures. In the first formulation (33), measures are supported on the set  $Q \times U$  of dimension 4+1=5 and we expect them to be located on the optimal trajectory  $(q^*(\cdot), u^*(\cdot))$ . An alternative proposed in [12] is to model controls by measures such that the measures are supported on Q only. Indeed, note that the dynamic in (32) is affine in the control u which takes its values inside the polytope  $conv\{-1,+1\}$ . 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. Following [12], this leads to consider

$$T_{LP'} = \inf_{\mu_1, \mu_2, \mu_f} \int d\mu_f$$

$$\int \left(\frac{\partial v}{\partial t}(t) + \frac{\partial v}{\partial q}(q)V(q)\right) d\mu = \int v(\cdot, q_f) d\mu_f - v(0, q_i), \quad \forall v \in \mathbb{R}[t, q], \quad (34)$$

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

where T is fixed,  $Q_f = [0, T]$ ,  $Q = B_2 \times B_2$  are admissible state sets.

#### • Step 2: Moment SDP.

An important feature of the problems (33)-(34) is their algebraic structure: the dynamic, the Lagrangian are polynomials and the sets Q, U are compact basic semialgebraic sets. In these settings, it is possible to handle the measures by their moments which yield a semi-definite program on countably many moments.

For a multi-index  $\alpha = (\alpha_1, \dots, \alpha_p) \in \mathbb{N}^p$ ,  $|\alpha|_1 = \sum_{i=1}^p \alpha_i$  and for  $y = (y_1, \dots, y_p)$ , the notation  $y^{\alpha}$  stands for the monomial  $y_1^{\alpha_1} \dots y_p^{\alpha_p}$ . We denote by  $\mathbb{N}_d^p$  the set  $\{\alpha \in \mathbb{N}^p \mid |\alpha|_1 \leq d\}$ .

Definition 2. The moment of order  $\alpha \in \mathbb{N}^p$  of a measure  $\mu$  supported on  $Z \subset \mathbb{R}^p$  is the real  $y_{\alpha} = \int z^{\alpha} d\mu$ .

 $\mu \in \mathcal{M}(z)$  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 3. Given an arbitrary sequence of real  $(y_{\alpha})_{\alpha}$ , we define the Riesz linear functional  $l_{\nu}: \mathbb{R}[z] \to \mathbb{R}$  by  $l_{\nu}(z^{\alpha}) = y_{\alpha}$  for all  $\alpha \in \mathbb{N}^{p}$ .

Definition 4. 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  $M_d(y)[i, j] = l_y(z^{i+j}) = y_{i+j}, \forall i, j \in \mathbb{N}_d^p$ .

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

**Proposition 12.** Let X be a compact basic semialgebraic set  $Z = \{z \in \mathbb{R}^p \mid g_k(z) \geq 0, k = 1, \dots, n_Z\}$ . Then, a necessary condition for a sequence  $(y_\alpha)_\alpha$  to have a representing measure  $\mu \in \mathcal{M}^+(z)$  is

$$M_d(y) \succeq 0$$
,  $M_d(g_k y) \succeq 0$ ,  $\forall d \in \mathbb{N}$ .

Write

$$[0,T] \times Q \times U = \{(t,q,u) \mid q = (q_{11},q_{12},q_{21},q_{22}), g_1(t,q,u) = t(T-t) \ge 0, g_i(t,q,u) = 1 - q_{i1}^2 - q_{i2}^2 \ge 0, i = 2,3, g_4(t,q,u) = 1 - u^2 \ge 0\},$$

 $Q_f = \{t \mid g^f(t) = t \, (T-t) \ge 0\}$  and denote by  $l_{y^\mu}, l_{y^{\mu_f}}$  the Riesz functionals associated respectively with two sequences  $y^\mu$  and  $y^{\mu_f}$ . Then, the moment SDP problem associated with (33) is

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

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

$$M_d(y^{\mu}) \succeq 0, M_d(g_i y^{\mu}) \succeq 0, \forall d \in \mathbb{N}, i = 1 \dots n_{Q \times U}$$

$$M_d(y^{\mu_f}) \succeq 0, M_d(g^f y^{\mu_f}) \succeq 0, \forall d \in \mathbb{N}.$$
(35)

According to Proposition 12, we have  $T_{LP} \ge T_{SDP}$  and this is in fact an equality according to [24, Theorem 3.8].

• Step 3: Hierarchy of SDP problems.

Note that  $M_{d+1}(y) \succeq 0$  implies  $M_d(y) \succeq 0$ . The LMI constraints and the sequence  $(y_\alpha)$  of (35) are truncated which yield for  $r \ge 1$  the Lasserre's hierarchy parameterized by  $r \ge 1$ 

$$T_{LMI}^{r} = \inf_{(y_{\alpha}^{\mu})_{|\alpha| \leq 2r}, (y_{\alpha}^{\mu_{f}})_{|\alpha| \leq 2r}} l_{y^{\mu_{f}}}(1)$$

$$l_{y^{\mu}} \left(\frac{\partial v}{\partial t} + \frac{\partial v}{\partial q}(q)V(q)\right) = l_{y^{\mu_{f}}}(v(\cdot, q_{f})) - v(0, q_{i}), \quad \forall v \in \mathbb{R}[t, q], \qquad (36)$$

$$M_{r}(y^{\mu}) \succeq 0, M_{r-s_{i}}(g_{i}y^{\mu}) \succeq 0, i = 1 \dots n_{Q \times U}$$

$$M_{r}(y^{\mu_{f}}) \succ 0, M_{r}(g^{f}y^{\mu_{f}}) \succ 0.$$

where  $s_i = \deg(g_i)/2$  if  $\deg(g_i)$  is even and  $s_i = (\deg(g_i) + 1)/2$  otherwise.

Proposition 13 (Theorem 5.6, [24]). We have

$$T_{SDP} \ge \ldots \ge T_{LMI}^{r+1} \ge T_{LMI}^r \ge \ldots \ge T_{LMI}^1$$
.

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

Numerical simulations for the multisaturation problem The problem (36) corresponds to the multisaturation problem of two spins associated with the LP problem (33). Likewise, we consider the LMI hierarchy associated with the LP problem (34) in the single spin case and the two spins case. We use the MOSEK toolbox [29] to solve the SDP problems. Let  $t_f$  denote the best solution found with the Bocop software and given in Table 1 for the single spin case or the two spins case. The value of the parameters for (36) are:  $\varepsilon = 0.1$ ,  $q_i = (0,1,0,1)$ ,  $q_f = (0,0,0,0)$  and  $T = t_f$ . In Fig. 16 are represented the relative error  $err(r) = |T_{LMI}^r - t_f|/t_f$  for the cases

 $C_1, C_2, C_3, C_4$  where  $T_{LMI}^r$  is the optimal value of (36) in the single spin case and the two spins case. The lower bounds of  $T_{min}^*$  are  $T_{LMI}^r$  (resp.  $T_{LMI'}^r$ ) associated with the LP problem (33) (resp. (34)) and are given in Table 2. The number of moments for the two problems associated with (33) and (34) are given by  $N_m = \binom{n+m+1+2d}{n+m+1} + \binom{n+1+2d}{n+1}$  for the 1st formulation and  $N_m = \binom{n+1+2d}{n+1}$  for the 2nd formulation where  $n = \dim Q$ ,  $m = \dim U$  and  $n_d = 2$  is the number of modes. 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.

	1st Fo	rmulation	2nd Formulation		
r	$N_m$	$T_{LMI}^r/t_f$	$N_m$	$T_{LMI'}^r/t_f$	
1	25	0.1857	30	0.182	
2	105	0.4836	105	0.4042	
3	294	0.7389	252	0.5645	
4	660	0.8509	495	0.8158	
5	1287	0.907	858	0.8716	
6	2275	0.936	1365	0.904	
7	3740	0.9483	2040	0.9203	
8	5814	0.9540	2907	0.9284	

Table 2: Single spin saturation for the case  $C_2$ .  $T_{LMI}^r$  is the optimal value of the hierarchy (36) for both formulations,  $t_f = 110.44$  is the time found by the Bocop software.

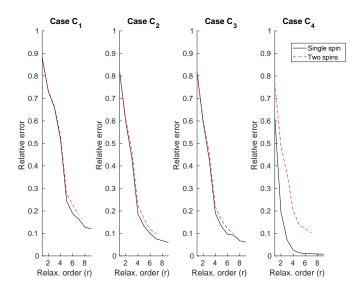


Figure 16: Saturation problem of one spin and two spins for the cases  $C_1, C_2, C_3, C_4$ . Relative error  $err(r) = |T_{LMI'}^r - t_f|/t_f$  where r is the order of relaxation,  $T_{LMI'}^r$  is the optimal value of (36) using the formulation (34) and  $t_f$  is the final time of (32) computed by Bocop.

#### 4.2.3 Homotopy with the indirect approach

Time-minimal saturation of a single spin. In the single-spin case, we illustrate the two behaviors described on Figure 4 on 3 examples, in Figures 17, 18 and 19. These figures represent the optimal trajectory and control for 3 spins in a family parameterized by  $\lambda$  as follows. Starting with the  $C_1$  case (deoxygenated blood) with relaxation times  $T_{1,0}$  and  $T_{2,0}$ , we consider the spin  $(T_1(\lambda), T_{2,0})$  with  $T_1(\lambda) = T_1 - d\lambda$ ,  $d = (2T_1 - 3T_2)/2$ .

In particular, with  $\lambda = 0$ , the spin is exactly deoxygenated blood. With  $\lambda = 1$ ,  $2\Gamma = 3\gamma$ . And with  $\lambda = \lambda_f = (T_1 - T_2)/d$ , the spin is isotrope ( $\delta = 0$ ) and it corresponds to the "water-like" case  $C_4$ .

As described in the introduction, we observe a bifurcation when the horizontal singular line at altitude  $z_S = \gamma/2\delta$  leaves the Bloch ball, that is when  $2\Gamma = 3\gamma$ . The experimental value of  $\lambda$  corresponding to the bifurcation is not exactly  $\lambda = 1$ , but rather a value  $\bar{\lambda} \simeq 1$  (here  $\bar{\lambda} \simeq 0.9941$ ). This is partially due to physical limitations on the control: the lowest point reachable by a trajectory is not the South pole of the Bloch ball.

Time-minimal saturation of two spins with  $B_1$ -inhomogeneity. In the multisaturation case, similar to the single spin case, bifurcations of BC-extremals (the optimality candidates given by the maximum principle) are described on Figs.20-21-22-23. The control was computed with homotopy method, using HamPath in combination with Bocop (in order to determine the structure of the extremal trajectory for  $\lambda=0$ ).  $\lambda=0$  corresponds to the  $C_1$  case and  $\lambda=\lambda_f$  corresponds to the  $C_4$  case. In particular, Fig.20 and Fig 23 have to be compared with the first and fourth graphs of Fig.15: there are many local solutions to the multisaturation problem with comparable costs, hence the crucial use of the LMI method. Figs.20-21 show a control with the same structure  $\gamma_+\gamma_s\gamma_+\gamma_s\gamma_+\gamma_s$ , that is three sequences of bang-singular. A bifurcation occurs at  $\bar{\lambda}\simeq 0.94$  where the first singular arc disappears. Figs.22-23 show a control with a different structure  $\gamma_+\gamma_s\gamma_+\gamma_s$ . In each figure, we have represented the critical altitude  $z_s=\gamma/2/\delta$  (represented with a horizontal dotted line on the figures). At  $\lambda=\lambda_f$  the extremal is simply  $\gamma_+\gamma_0\gamma_+\gamma_0$ : singular arcs are obtained by applying a zero control.

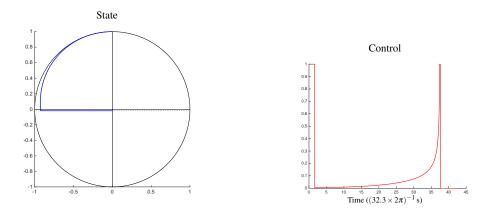


Figure 17: Time-minimal saturation of a single spin ( $\lambda = 0$ )

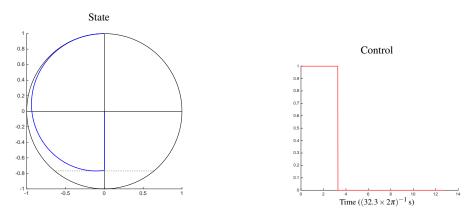


Figure 18: Time-minimal saturation of a single spin  $(\lambda=0.9941\simeq\bar{\lambda})$ 

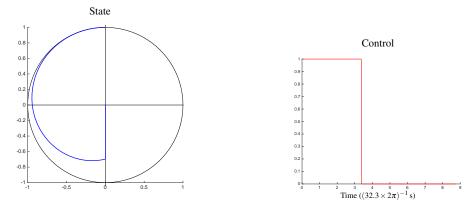


Figure 19: Time-minimal saturation of a single spin  $(\lambda = \lambda_f)$ 

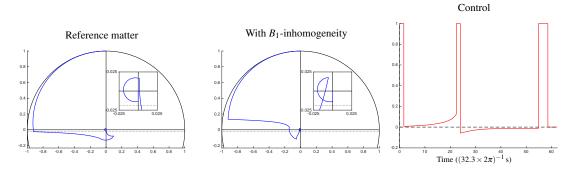


Figure 20: BC-extremal for the multisaturation problem with  $\lambda=0$ 

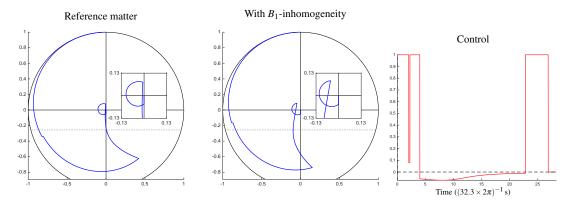


Figure 21: BC-extremal for the multisaturation problem with  $\lambda=0.943<\bar{\lambda}$ 

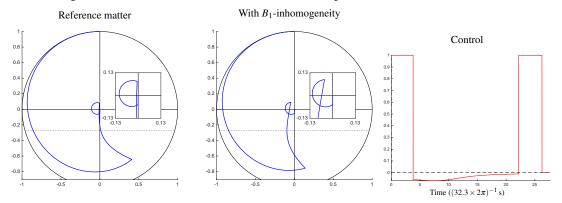


Figure 22: BC-extremal for the multisaturation problem with  $\lambda=0.948>\bar{\lambda}$ 

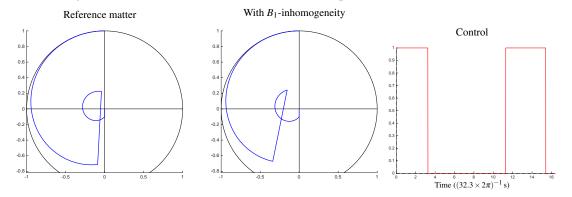


Figure 23: BC-extremal for the multisaturation problem with  $\lambda=\lambda_f$ 

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