

A convex analysis approach to multiswitching control of partial differential equations

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A convex analysis approach for computing multiswitching controls for partial differential equations is introduced; such controls consist of an arbitrary number of components of which at most one should be simultaneously active. Using a Moreau–Yosida approximation, a family of approximating problems is obtained that is amenable to solution by a semismooth Newton method. The efficiency of this approach and the structure of the obtained controls are demonstrated by numerical examples.

1 Introduction

Multiswitching control refers to time-dependent optimal control problems with several control mechanisms of which at most one should be active at every point in time. We shall consider here optimal tracking control for a linear diffusion equation $Ly = Bu$ on $\Omega_T := (0, T] \times \Omega$, $y(0) = y_0$ on Ω , where $L = \partial_t - A$ for an elliptic operator A defined on $\Omega \subset \mathbb{R}^n$ carrying suitable boundary conditions. The control operator B is defined by

$$(Bu)(t, x) = \sum_{i=1}^N \chi_{\omega_i}(x) u_i(t),$$

where χ_{ω_i} is the characteristic function of the given control domain $\omega_i \subset \Omega$ of positive measure. Furthermore, let $\omega_{\text{obs}} \subset \Omega_T$ denote the observation domain and let $y^d \in L^2(\omega_{\text{obs}})$ denote the

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target. We aim to promote the switching structure of the optimal control $\bar{u} \in L^2(0, T; \mathbb{R}^N)$ by a convex penalty term and consider for this purpose the optimal control problem

$$(P) \quad \begin{cases} \min_{u \in L^2(0, T; \mathbb{R}^N)} \frac{1}{2} \|y - y^d\|_{L^2(\omega_{\text{obs}})}^2 + \frac{\alpha}{2} \int_0^T |u(t)|_1^2 dt, \\ \text{s. t. } Ly = Bu, \quad y(0) = y_0. \end{cases}$$

This is a convex optimization problem, for which we derive first-order optimality conditions in primal-dual form whose Moreau–Yosida regularization (to be introduced below) can be solved using a superlinearly convergent semismooth Newton method.

The approach which we follow here is related to both the switching control problem in [5] and the distributed parabolic sparse control problem in [4]. In the former, a nonconvex formulation in the case where $N = 2$ was considered; we compare its convex relaxation to the present approach below. In the latter, the L^2 norm in time of the measure norm in space was used to promote temporally varying sparsity in space. While the choice of the nonsmooth functional involving the controls in (P) is motivated by sparsity considerations, one can arrive at this functional as well from controllability–observability considerations. In fact, it was shown in [19, Theorem 4.1] that – provided an appropriately defined controllability Gramian has full rank – exact null controls with perfect switching have minimal $L^2(0, T; \mathbb{R}^2)$ norm, where \mathbb{R}^2 is endowed with the ℓ^1 norm.

Let us comment on further related work. While there is a large body of work on the stability of switching control for ordinary differential equations (see, e.g., [3, 15]), the corresponding problem for time-dependent partial differential equations is less well-studied. In both cases, one should distinguish switching control in the sense defined above from the control of switched systems, where a binary or integer decision variable is sought that toggles between different forms of the state equation. For the latter in the context of PDEs, we refer to [8, 16]. Time-dependent switching controls can be realized in this framework by considering a product of a binary decision variable and a continuous control, which leads to a bilinear mixed-integer optimal control problem, see [9]. In contrast, we aim directly for continuous controls with the desired switching structure without introducing auxiliary decision variables. Exact null controls with such switching structure for the heat and wave equation were treated in [19, 14, 18] and [7], respectively. In addition, most of these works restrict their analysis to the case of switching between two control components. Here, in contrast, we treat optimal control problems with multiple switching control components.

Our work is organized as follows. Section 2 discusses the existence and first-order optimality conditions for solutions to (P) as well as its regularization within an abstract convex analysis framework. Explicit pointwise characterizations of the switching relations arising from the optimality system and for its regularization are given in Section 3. Here we also discuss the relation of the proposed switching functional in (P) to other possible choices of the penalty term. Section 4 is concerned with the numerical solution of the regularized optimality system by a semismooth Newton method. A numerical example for multiswitching control of a two-dimensional linear heat equation is computed in Section 5.

2 Convex analysis approach

We recall the convex analysis approach for (nonconvex) switching controls for partial differential equations from [5], which is also applicable to the convex penalty considered here. For this purpose, we consider Problem (P) in the reduced form

$$\min_u \mathcal{F}(u) + \mathcal{G}(u),$$

with $\mathcal{F} : L^2(0, T; \mathbb{R}^N) \rightarrow \mathbb{R}$ and $\mathcal{G} : L^2(0, T; \mathbb{R}^N) \rightarrow \mathbb{R}$ given by

$$\mathcal{F}(u) = \frac{1}{2} \|Su - y^d\|_{L^2(\omega_{\text{obs}})}^2, \quad \mathcal{G}(u) = \frac{\alpha}{2} \int_0^T |u(t)|_1^2 dt,$$

where the continuous affine solution operator $S : u \mapsto y$ assigns to any control $u \in L^2(0, T; \mathbb{R}^N)$ the unique state $y \in L^2(\Omega_T)$ satisfying the state equation $Ly = Bu$ with initial condition $y(0) = y_0$ subject to appropriate boundary conditions. Here we assume that the coefficients of A , the boundary and initial conditions as well as the domain Ω are sufficiently regular that the range of S is contained in

$$(2.1) \quad W(0, T) := L^2(0, T; H_0^1(\Omega)) \cap W^{1,2}(0, T; H^{-1}(\Omega)) \hookrightarrow C([0, T]; L^2(\Omega)).$$

Since S is affine, \mathcal{F} is proper, convex and lower semicontinuous. Furthermore, since the squared norm $|\cdot|_1^2$ is convex, \mathcal{G} is proper, convex, lower semicontinuous, and, in addition, radially unbounded. Existence of a solution thus follows from standard arguments, e.g., Tonelli's direct method.

Proposition 2.1. *There exists a minimizer \bar{u} for Problem (P).*

We next derive first-order optimality conditions in primal-dual form. Throughout, for any proper convex function \mathcal{H} , we denote by \mathcal{H}^* its Fenchel conjugate and by $\partial\mathcal{H}$ its subdifferential; see, e.g., [1, 6] for their definitions. The following proposition is a direct consequence of the sum rule and inversion formula for convex subdifferentials (see, e.g., [1, Corollary 16.24] for the latter) as well as the Fréchet-differentiability of \mathcal{F} .

Proposition 2.2. *The control $\bar{u} \in L^2(0, T; \mathbb{R}^N)$ is a minimizer for (P) if and only if there exists a $\bar{p} \in L^2(0, T; \mathbb{R}^N)$ such that*

$$(OS) \quad \begin{cases} -\bar{p} = \mathcal{F}'(\bar{u}), \\ \bar{u} \in \partial\mathcal{G}^*(\bar{p}), \end{cases}$$

holds.

Since \mathcal{F} is a standard quadratic tracking term, the first relation in (OS) can be expressed in a straightforward manner in terms of the solution operator $S = L^{-1}B$ and its adjoint $S^* = B^*L^{-*}$ (with homogeneous boundary and initial conditions), i.e., $\bar{p} = -S^*(S\bar{u} - y^d)$. For later use, we point out that due to (2.1) and the specific choice of S there holds $\bar{p} \in V := B^*(W(0, T)) \hookrightarrow L^r(0, T; \mathbb{R}^N)$ for any $r > 2$.

The second relation is responsible for the switching structure of the optimal control \bar{u} , and we will give a pointwise characterization in Proposition 3.1 below.

Our numerical approach is based on the Moreau–Yosida regularization of (OS). Specifically, we replace $\partial\mathcal{G}^*$ for $\gamma > 0$ by

$$\partial\mathcal{G}_\gamma^*(p) := (\partial\mathcal{G}^*)_\gamma(p) := \frac{1}{\gamma} \left(p - \text{prox}_{\gamma\mathcal{G}^*}(p) \right),$$

where

$$\text{prox}_{\gamma\mathcal{G}^*}(v) := \arg \min_w \frac{1}{2\gamma} \|w - v\|^2 + \mathcal{G}^*(w) = (\text{Id} + \gamma\partial\mathcal{G}^*)^{-1}(v)$$

is the proximal mapping of \mathcal{G}^* , which in Hilbert spaces coincides with the resolvent of $\partial\mathcal{G}^*$. Note that the proximal mapping and thus the Moreau–Yosida regularization of a proper and convex functional is always single-valued and Lipschitz continuous; see, e.g., [1, Corollary 23.10].

We then consider the regularized system

$$(OS_\gamma) \quad \begin{cases} -p_\gamma = \mathcal{F}'(u_\gamma), \\ u_\gamma = \partial\mathcal{G}_\gamma^*(p_\gamma). \end{cases}$$

Again, we will give an explicit formulation of (OS_γ) in the next section.

Proposition 2.3. *For each $\gamma > 0$ system (OS_γ) admits a unique solution (u_γ, p_γ) .*

Proof. Using convex analysis techniques (see, e.g., [1, Chapter 12]), we obtain that (OS_γ) is the necessary optimality condition for

$$(P_\gamma) \quad \min_u \mathcal{F}(u) + \left(\mathcal{G}_\gamma^*\right)^*(u).$$

In fact, since \mathcal{F} is globally defined and continuous, we have the necessary optimality condition

$$0 \in \partial\mathcal{F}(u_\gamma) + \partial\left(\mathcal{G}_\gamma^*\right)^*(u_\gamma).$$

Setting $p_\gamma := -\mathcal{F}'(u_\gamma)$ we recall again the subdifferential inversion formula from [1, Corollary 16.24]. Additionally, we note that the Yosida-regularization $\partial\mathcal{G}_\gamma^*$ of the subdifferential $\partial\mathcal{G}^*$ coincides with the Fréchet derivative of the Moreau-envelope \mathcal{G}_γ^* of \mathcal{G}^* ; see, e.g., [1, Proposition 12.29]. Together, these yield (OS_γ) .

We turn now to the existence and uniqueness of a solution to (OS_γ) . From [1, Remark 12.24], we have the alternative characterization of the Moreau-envelope via the infimal convolution

$$\mathcal{G}_\gamma^* = \mathcal{G}^* \square \frac{1}{2\gamma} \|\cdot\|_{L^2(0,T;\mathbb{R}^N)}^2,$$

where \mathbb{R}^N is endowed with the Euclidean norm. Furthermore, from [1, Theorem 15.3], we obtain

$$\begin{aligned} \left(\mathcal{G}_\gamma^*\right)^* &= \left(\mathcal{G}^* \square \frac{1}{2\gamma} \|\cdot\|_{L^2(0,T;\mathbb{R}^N)}^2\right)^* = \left(\left(\mathcal{G} + \frac{\gamma}{2} \|\cdot\|_{L^2(0,T;\mathbb{R}^N)}^2\right)^*\right)^* \\ &= \mathcal{G} + \frac{\gamma}{2} \|\cdot\|_{L^2(0,T;\mathbb{R}^N)}^2. \end{aligned}$$

This implies that $(\mathcal{G}_\gamma^*)^*$ is strictly convex. Moreover, $\mathcal{G}_\gamma^* \leq \mathcal{G}^*$ and hence $0 \leq \mathcal{G}^{**} = \mathcal{G} \leq (\mathcal{G}_\gamma^*)^*$. Therefore a minimizing sequence for (P_γ) is necessarily bounded, and by a weak subsequential limit argument, existence of a unique solution u_γ to (P_γ) and thus of $p_\gamma = -\mathcal{F}'(u_\gamma)$ satisfying (OS_γ) follows. \square

Finally, convergence as $\gamma \rightarrow 0$ can be shown as in [5, Proposition 2.5]. This requires showing that $\{\partial\mathcal{G}^*(p_\gamma)\}_{\gamma>0}$ is uniformly bounded in γ provided that $\{p_\gamma\}_{\gamma>0}$ is bounded. But this follows from the explicit characterization in (3.3) below.

Proposition 2.4. *For any sequence $\gamma_n \rightarrow 0$, the sequence $\{(u_{\gamma_n}, p_{\gamma_n})\}_{n \in \mathbb{N}}$ converges weakly to a solution (\bar{u}, \bar{p}) to (OS) .*

3 Multiswitching penalty

Here we compute an explicit, pointwise, characterization of $\partial\mathcal{G}^*$ and $\partial\mathcal{G}_\gamma^*$ by exploiting the fact that for integral functionals of the form

$$\mathcal{G}(u) = \int_0^T g(u(t)) dt,$$

the Fenchel conjugate and convex subdifferential can be computed pointwise via g ; see, e.g., [6, Props. IV.1.2, IX.2.1], [1, Prop. 16.50]. We thus focus on

$$g : \mathbb{R}^N \rightarrow \mathbb{R}, \quad g(v) = \frac{\alpha}{2} |v|_1^2.$$

Because the Fenchel conjugate of half the squared norm is half the squared dual norm (see, e.g., [2, Example 3.27]), a scaling argument (e.g., [1, Proposition 13.20 (i)]) yields

$$(3.1) \quad g^* : \mathbb{R}^N \rightarrow \mathbb{R}, \quad g^*(q) = \frac{1}{2\alpha} |q|_\infty^2 = \max_{1 \leq i \leq N} \frac{1}{2\alpha} q_i^2.$$

3.1 Subdifferential

Since g^* is the maximum of finitely many convex and differentiable functions, its convex subdifferential is given by

$$\partial g^*(q) = \overline{\text{co}} \left(\bigcup_{\{i: g^*(q) = g_i^*(q)\}} \{(g_i^*)'(q)\} \right),$$

where $\overline{\text{co}}$ denotes the closed convex hull and $g_i^*(q) = \frac{1}{2\alpha} q_i^2$; see, e.g., [10, Corollary 4.3.2]. It is instructive to first consider the case $N = 2$. In this case,

$$(3.2) \quad \partial g^*(q) = \begin{cases} \left\{ \left(\frac{1}{\alpha} q_1, 0 \right) \right\} & \text{if } |q_1| > |q_2|, \\ \left\{ \left(0, \frac{1}{\alpha} q_2 \right) \right\} & \text{if } |q_1| < |q_2|, \\ \left\{ t \left(\frac{1}{\alpha} q_1, 0 \right) + (1-t) \left(0, \frac{1}{\alpha} q_2 \right) : t \in [0, 1] \right\} & \text{if } |q_1| = |q_2|. \end{cases}$$

For the general case, the subdifferential can be given componentwise for $1 \leq j \leq N$ as

$$(3.3) \quad [\partial g^*(q)]_j = \begin{cases} \left\{ \frac{1}{\alpha} q_j \right\} & \text{if } |q_j| = \max_i |q_i| \text{ and } |q_j| > |q_i|, i \neq j, \\ \{0\} & \text{if } |q_j| < \max_i |q_i|, \\ \left\{ \frac{s_j}{\alpha} q_j : s_j \geq 0, \sum s_i = 1 \right\} & \text{if } |q_j| = \max_i |q_i| \text{ and } |q_j| = |q_i| \text{ for } i \neq j, \end{cases}$$

where the sum in the last set is over all i with $|q_i| = |q_j|$. We thus obtain for $u, p \in L^2(0, T; \mathbb{R}^N)$ the pointwise characterization

$$\partial \mathcal{G}^*(p) = \left\{ u \in L^2(0, T; \mathbb{R}^N) : u(t) \in \partial g^*(p(t)) \text{ for a.e. } t \in (0, T) \right\}.$$

In particular, this yields a pointwise characterization of the second relation in (OS).

Proposition 3.1. *The minimizer $\bar{u} \in L^2(0, T; \mathbb{R}^N)$ of (P) and $\bar{p} := -\mathcal{F}'(\bar{u}) \in L^2(0, T; \mathbb{R}^N)$ satisfy for almost every $t \in (0, T)$ and $1 \leq j \leq N$*

$$\bar{u}_j(t) \in \begin{cases} \left\{ \frac{1}{\alpha} \bar{p}_j(t) \right\} & \text{if } |\bar{p}_j(t)| = \max_i |\bar{p}_i(t)| \text{ and } |\bar{p}_j(t)| > |\bar{p}_i(t)|, i \neq j, \\ \{0\} & \text{if } |\bar{p}_j(t)| < \max_i |\bar{p}_i(t)|, \\ \left\{ \frac{s_j}{\alpha} \bar{p}_j(t) : s_j \geq 0, \sum s_i = 1 \right\} & \text{if } |\bar{p}_j(t)| = \max_i |\bar{p}_i(t)| \text{ and } |\bar{p}_j(t)| = |\bar{p}_i(t)| \text{ for } i \neq j. \end{cases}$$

From this proposition, the desired switching property follows: Outside of the singular arc

$$\mathcal{S} := \left\{ t \in (0, T) : |\bar{p}_{j_1}(t)| = |\bar{p}_{j_2}(t)| = \max_i |\bar{p}_i(t)| \text{ for } j_1 \neq j_2 \right\},$$

we have $\bar{u}_{j_1} \bar{u}_{j_2} = 0$ for all $j_1 \neq j_2$. In particular, if \mathcal{S} has Lebesgue measure zero, \bar{u} exhibits perfect switching, i.e., $\bar{u}_{j_1} \bar{u}_{j_2} = 0$ almost everywhere. Under the stronger assumption that $|\bar{p}_{j_1}(t)| \neq |\bar{p}_{j_2}(t)|$ for all $j_1 \neq j_2$ and almost all $t \in (0, T)$, a similar relation for exact null controls was given in [19, equations (2.36–2.38)].

Remark 1. *It is difficult to give general conditions for the optimal control to be perfectly switching; instead one would have to exploit properties of the specific (adjoint) state equation. For example, in the case of the one-dimensional heat equation, one could use the fact that solutions are real-analytic with respect to time to argue that $|p_i(t)| = |p_j(t)|$ for all $t \in O$ for an open set $O \subset (0, T)$ is only possible if $p_i(t) = p_j(t) = 0$; cf. [19, Section 6.1].*

3.2 Proximal mapping

We next characterize the proximal mapping $\text{prox}_{\gamma g^*}(p)$ via the resolvent $(\text{Id} + \gamma \partial g^*)^{-1}(p)$ by proceeding similarly to [5, Section 3.3]; see also [12].

Proposition 3.2. *Let the components of $v \in \mathbb{R}^N$ be sorted by decreasing magnitude and let d be the smallest index for which*

$$(3.4) \quad |v_{d+1}| < \frac{\alpha}{d\alpha + \gamma} \sum_{i=1}^d |v_i|$$

holds; if no such index exists, let $d = N$. The explicit form of the proximal mapping is then given componentwise for $1 \leq j \leq N$ by

$$[\text{prox}_{\gamma g^*}(v)]_j = \begin{cases} \text{sign}(v_j) \frac{\alpha}{d\alpha + \gamma} \sum_{i=1}^d |v_i| & \text{if } j \leq d, \\ v_j & \text{if } j > d. \end{cases}$$

Proof. Let $w := (\text{Id} + \gamma \partial g^*)^{-1}(v)$. This is characterized by the subdifferential inclusion

$$(3.5) \quad v \in (\text{Id} + \gamma \partial g^*)(w) = \{w\} + \gamma \partial g^*(w).$$

First, we observe that (3.5) implies that $\text{sign}(v_i) = \text{sign}(w_i)$. We next follow the case discrimination in the characterization of the subdifferential, where we assume that the components of w are ordered by magnitude, i.e., $|w_1| \geq |w_2| \geq \dots \geq |w_N|$ (the remaining cases following by permutation).

- (i) $|w_1| > |w_2|$: In this case, all components of the subdifferentials are single-valued; solving for each w_i in (3.5) yields

$$w_i = \begin{cases} \frac{\alpha}{\alpha + \gamma} v_i & \text{if } i = 1, \\ v_i & \text{if } i > 1. \end{cases}$$

The assumption $|w_1| > |w_2|$ is then equivalent to the condition

$$|v_1| > \left(1 + \frac{\gamma}{\alpha}\right) |v_2|.$$

- (ii) $|w_1| = |w_2| > |w_3|$: Using that $\text{sign}(w_i) = \text{sign}(v_i)$, this implies that $w_1 = \text{sign}(v_1)\bar{w}$ and $w_2 = \text{sign}(v_2)\bar{w}$ for some $\bar{w} > 0$. Inserting this into the set-valued case of the subdifferential, we deduce that for some $t \in [0, 1]$,

$$\begin{aligned} \text{sign}(v_1)|v_1| &= \left(1 + t \frac{\gamma}{\alpha}\right) \text{sign}(v_1)\bar{w}, \\ \text{sign}(v_2)|v_2| &= \left(1 + (1-t) \frac{\gamma}{\alpha}\right) \text{sign}(v_2)\bar{w}. \end{aligned}$$

Dividing by the $\text{sign}(v_i)$ on both sides and adding yields

$$|v_1| + |v_2| = \left(2 + \frac{\gamma}{\alpha}\right) \bar{w},$$

from which we obtain

$$w_i = \begin{cases} \text{sign}(v_i) \frac{\alpha}{2\alpha + \gamma} (|v_1| + |v_2|), & \text{if } i \leq 2, \\ v_i & \text{if } i > 2, \end{cases}$$

the single-valued cases for $i > 2$ following as before.

Inserting these values into $|w_2| > |w_3|$ then yields the condition

$$|v_1| + |v_2| > \left(2 + \frac{\gamma}{\alpha}\right) |v_3|.$$

Note that in addition,

$$|v_1| \leq \left(1 + \frac{\gamma}{\alpha}\right) |v_2|$$

must hold since due to case (i), the converse is equivalent to $|w_1| > |w_2|$, in contradiction to the assumption.

(iii) $|w_1| = |w_2| = |w_3| > |w_4|$: As before, the chain of equalities is equivalent to

$$\begin{aligned} |v_1| &= \left(1 + t_1 \frac{\gamma}{\alpha}\right) \bar{w}, \\ |v_2| &= \left(1 + t_2 \frac{\gamma}{\alpha}\right) \bar{w}, \\ |v_3| &= \left(1 + (1 - t_1 - t_2) \frac{\gamma}{\alpha}\right) \bar{w} \end{aligned}$$

for some t_1, t_2 with $t_1 + t_2 \in (0, 1)$. Adding these equations yields that

$$w_i = \begin{cases} \text{sign}(v_i) \frac{\alpha}{3\alpha + \gamma} (|v_1| + |v_2| + |v_3|), & \text{if } i \leq 3, \\ v_i & \text{if } i > 3, \end{cases}$$

as well as the condition

$$|v_1| + |v_2| + |v_3| > \left(3 + \frac{\gamma}{\alpha}\right) |v_4|.$$

Again,

$$|v_1| \leq \left(1 + \frac{\gamma}{\alpha}\right) |v_2|, \quad \text{and} \quad |v_1| + |v_2| \leq \left(2 + \frac{\gamma}{\alpha}\right) |v_3|$$

must hold since otherwise we obtain a contradiction in the case distinction.

(iv) The remaining cases $|w_1| = \dots = |w_j| > |w_{j+1}|$ for $j = 4, \dots, N$ follow analogously, establishing the desired characterization of the proximal mapping. \square

It is now straight-forward to give a pointwise characterization of the Moreau–Yosida regularization $\partial \mathcal{G}_\gamma^*$ via

$$[\partial \mathcal{G}_\gamma^*(p)](t) = \partial g_\gamma^*(p(t)) = \frac{1}{\gamma} \left(p(t) - \text{prox}_{\gamma g^*}(p(t)) \right) \quad \text{for a.e. } t \in (0, T).$$

We point out that for every $t \in (0, T)$, the number d in Proposition 3.2 is the number of nonzero components in $u_\gamma(t) = [\partial \mathcal{G}_\gamma^*(p)](t)$. In particular for the case $d = 1$, the explicit form of ∂g_γ^* is given componentwise as

$$[\partial g_\gamma^*(q)]_j = \begin{cases} \frac{1}{\alpha + \gamma} q_j & \text{if } |q_j| = \max_i |q_i| \text{ and } |q_j| > \left(1 + \frac{\gamma}{\alpha}\right) |q_i|, i \neq j, \\ 0 & \text{if } |q_j| < \max_i |q_i|. \end{cases}$$

Thus if for some $\gamma > 0$, the solution (u_γ, p_γ) to (OS_γ) is such that $d = 1$ for almost every $t \in (0, T)$, the regularized control is perfectly switching. In this sense, the Moreau–Yosida regularization can lead to an exact penalization of the switching constraint, which is demonstrated in the numerical examples below.

3.3 Relation to other approaches for switching control

In this section, we address the relation of the proposed approach to the ℓ^0 formulation of [5] as well as to alternative ℓ^1 -type penalizations. For this purpose, we first point out that an equivalent formulation of the control cost term $g(v)$ is

$$(3.6) \quad g(v) = \frac{\alpha}{2}|v|_2^2 + \alpha \sum_{\substack{i,j=1 \\ i < j}}^N |v_i v_j|,$$

which includes an ℓ^1 penalization of $v_i v_j \neq 0$ for all $i \neq j$. We now compare this formulation with other approaches in the case of $N = 2$.

First, consider the nonconvex ℓ^0 -switching formulation from [5],

$$(3.7) \quad g_0(v) = \frac{\alpha}{2}|v|_2^2 + \beta|v_1 v_2|_0$$

(where $|s|_0 = 0$ if $s = 0$ and 1 otherwise), whose Fenchel conjugate is given by

$$(3.8) \quad g_0^*(q) = \begin{cases} \frac{1}{2\alpha} q_1^2 & \text{if } |q_1| \geq |q_2| \text{ and } |q_2| \leq \sqrt{2\alpha\beta}, \\ \frac{1}{2\alpha} q_2^2 & \text{if } |q_1| \leq |q_2| \text{ and } |q_1| \leq \sqrt{2\alpha\beta}, \\ \frac{1}{2\alpha}(q_1^2 + q_2^2) - \beta & \text{if } |q_1|, |q_2| \geq \sqrt{2\alpha\beta}. \end{cases}$$

For fixed q and β sufficiently large, this coincides with (3.2), since the third case (corresponding to the free arc, where both control components are active) is never attained. Therefore, the functional g can be interpreted as the convex relaxation of g_0 in the limit as $\beta \rightarrow \infty$, in the following sense: Taking the formal limit $\beta \rightarrow \infty$ in (3.7), we arrive at the (still nonconvex) constrained functional

$$g_\infty(v) := \frac{\alpha}{2}|v|_2^2 + \delta_{\{v: v_1 v_2 = 0\}}(v) := \begin{cases} \frac{\alpha}{2}|v|_2^2 & \text{if } v_1 v_2 = 0, \\ \infty & \text{else,} \end{cases}$$

where the switching property $v_i v_j = 0$ for all $i \neq j$ is included as an explicit constraint. Proceeding as in [5, § 3.2], we find that the Fenchel conjugate of g_∞ is given by the first two cases of (3.8). We thus have that $g_\infty^* = g^*$, from which it follows that

$$g = g^{**} = (g^*)^* = (g_\infty^*)^* = g_\infty^{**},$$

i.e., g is the lower convex envelope of g_∞ .

It would also be possible to have different penalty parameters for each term in (3.6), i.e., to consider for $\alpha, \beta > 0$

$$g_\beta(v) := \frac{\alpha}{2}|v|_2^2 + \beta|v_1 v_2|.$$

If $\beta < \alpha$, we can introduce $\gamma := \alpha - \beta > 0$ and rewrite g_β as

$$g_\beta(v) = \frac{\beta}{2}|v|_2^2 + \beta|v_1 v_2| + \frac{\gamma}{2}|v|_2^2 = (g_\gamma^*)^*(v).$$

The optimality system in this case is therefore equivalent to (OS_γ) .

In case $\beta > \alpha$, we can reformulate again to

$$g_\beta(v) = \frac{\alpha - \beta}{2} |v|_2^2 + \frac{\beta}{2} |v|_1^2,$$

which is obviously nonconvex. Computing the Fenchel conjugate leads to

$$g_\beta^*(q) = \begin{cases} \frac{1}{2\alpha} q_1^2 & \text{if } |q_1| \geq |q_2|, \\ \frac{1}{2\alpha} q_2^2 & \text{if } |q_1| \leq |q_2|, \end{cases}$$

which coincides with (3.1). As in the case of g_∞ , this implies that g is the convex envelope of g_β .

We conclude that our choice of g is the natural convex formulation for promoting switching controls.

4 Numerical solution

We apply a semismooth Newton method (see, e.g., [12, 17]) to (OS_γ) . By eliminating u_γ via the second relation of (OS_γ) and introducing the state $y_\gamma := S(u_\gamma) \in Y := L^2(\Omega_T)$, we arrive at the equivalent system

$$(4.1) \quad \begin{cases} y_\gamma = SH_\gamma(p_\gamma), \\ p_\gamma = -S^*(y_\gamma - y^d), \end{cases}$$

where we set $H_\gamma := \partial \mathcal{G}_\gamma^*$. We similarly set $h_\gamma := \partial g_\gamma^*$ such that

$$[H_\gamma(p)](t) = h_\gamma(p(t)) \quad \text{for a.e. } t \in (0, T).$$

Exactly as in [5, §5], one argues that $h_\gamma : \mathbb{R}^N \rightarrow \mathbb{R}^N$ is semismooth, and a Newton derivative $D_N h_\gamma(q) \in \mathbb{R}^{N \times N}$ is obtained by choosing any element of the Clarke derivative $\partial_C h_\gamma(q)$ (which for the piecewise differentiable function h_γ is given by the convex hull of the piecewise derivatives). Here we take, assuming again for simplicity that the components of q are sorted descending by magnitude, componentwise for $1 \leq i, j \leq N$

$$[D_N h_\gamma(q)]_{ji} = \begin{cases} \frac{(d-1)\alpha + \gamma}{\gamma(d\alpha + \gamma)} & \text{if } j = i \leq d, \\ -\frac{\alpha}{\gamma(d\alpha + \gamma)} \text{sign}(q_j q_i) & \text{if } j \leq d, i \leq d, i \neq j, \\ 0 & \text{if } j > d \text{ or } i > d, \end{cases}$$

where d is as in Proposition 3.2. The Newton derivative for arbitrary $q \in \mathbb{R}^N$ can be obtained from this by appropriate permutation of rows and columns. Note that $D_N h_\gamma(q)$ is symmetric. Furthermore, we point out that evaluation of the proximal mapping amounts to sorting for each $t \in (0, T)$ the vector $p(t) \in \mathbb{R}^N$, which can be carried out in $\mathcal{O}(N \log N)$ operations, and testing the $N - 1$ conditions in (3.4) for the sorted vector, which requires $\mathcal{O}(N)$ operations. Hence, evaluating h_γ and assembling the Newton derivative $D_N h_\gamma$ can be performed in $\mathcal{O}(N \log N)$ operations, thus avoiding the exponential complexity involved in computing switching points between any two (or more) control components; see, e.g., [11].

Since h_Y is semismooth from \mathbb{R}^N to \mathbb{R}^N , the corresponding superposition operator H_Y is semismooth from $V \hookrightarrow L^r(0, T; \mathbb{R}^N)$ for $r > 2$ to $L^2(0, T; \mathbb{R}^N)$, with Newton derivative at p in direction δp given by

$$[D_N H_Y(p) \delta p](t) = D_N h_Y(p(t)) \delta p(t) \quad \text{for a.e. } t \in (0, 1).$$

Considering the system (4.1) as an operator equation from $Y \times V$ to $Y \times V$, a semismooth Newton step for its solution thus consists of computing $(\delta y, \delta p) \in Y \times V$ for given $(y^k, p^k) \in Y \times V$ such that

$$(4.2) \quad \begin{cases} \delta y - S_0 D_N H_Y(p^k) \delta p = -y^k + S H_Y(p^k), \\ \delta p + S^* \delta y = -p^k - S^*(y^k - y^d), \end{cases}$$

and setting $y^{k+1} = y^k + \delta y$ and $p^{k+1} = p^k + \delta p$. Here, S_0 denotes the solution operator of the state equation with homogeneous initial and boundary conditions. As in [5, Proposition 5.2], one shows uniformly bounded invertibility of each Newton step, from which locally superlinear convergence of the semismooth Newton method follows.

The disadvantage of (4.2) is the forward–backward structure in time, which prevents using time-stepping methods for the application of S_0 and S^* . For the practical implementation, we therefore eliminate δy via the first relation of (4.2) as well as y^k via the first relation of (4.1) to obtain the reduced semismooth Newton step

$$(4.3) \quad \delta p + S^* S_0 D_N H_Y(p^k) \delta p = -\left(p^k + S^*(S H_Y(p^k) - y^d)\right),$$

which allows for time-stepping and matrix-free Krylov methods. The linear operator on the left-hand side becomes self-adjoint using the inner product generated by $D_N H_Y(p^k)$, and hence we apply a matrix-free CG method for the solution of (4.3) using this inner product; see [13]. We point out that monitoring the convergence of the CG method is still based on the standard Euclidean norm of the residual.

An alternative to (4.3) is a reduction to δy , which would lead to a reduced Newton matrix that is already self-adjoint with respect to the standard inner product. We do not follow this approach, because it would lead to a significantly higher dimension of the Newton system (full state space versus control space). In addition, the iterates would be infeasible, since in this case the state equation is only satisfied for a solution of the reduced optimality system.

To account for the local convergence of Newton methods, we embed the semismooth Newton method within a homotopy strategy for γ , where we start with a large γ which is successively reduced, taking the previous solution as starting point. Furthermore, we include a backtracking line search based on the residual norm to improve robustness.

5 Numerical examples

We now illustrate the behavior of the proposed approach and the structure of the resulting controls using a family of numerical examples. We consider the heat equation in two dimensions, i.e., $y = S(u)$ solves

$$y_t - \Delta y = \sum_{i=1}^N \chi_{\omega_i}(x) u_i(t) \quad \text{on } \Omega_T,$$

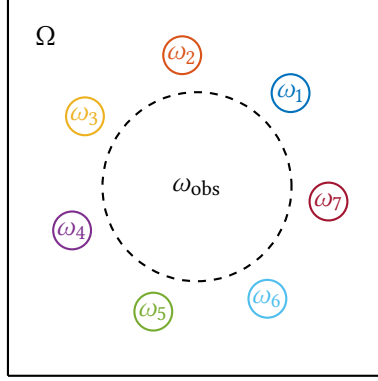


Figure 1: Problem setting for $N = 7$ control components

with $T = 10$ and $\Omega = (-1, 1)^2$, and zero initial and Neumann boundary conditions. We choose circular control domains $\omega_i = B_{0.1}(x_i)$ with centers $x_i = (\cos(\varphi_i), \sin(\varphi_i))/\sqrt{2}$ regularly distributed along a circle, where $\varphi_i = \pi/4 + 2\pi(i-1)/N$, while the observation domain is $\omega_{\text{obs}} = B_{0.5}(0)$; see Figure 1. The desired state $y^d \in L^2(\omega_{\text{obs}})$ is set to

$$y^d = \sum_{i=1}^N \cos(i+t) \sin^2\left(2\pi\frac{t}{T}\right) |x - x_i|^2.$$

The corresponding optimal controls are computed with the semismooth Newton method (4.3), stopping at a relative tolerance of 10^{-6} in the residual $F_k = \|F(p^k)\|_{L^2(0,T;\mathbb{R}^N)}$, where $F(p)$ denotes the right-hand side of (4.3). The backtracking line search starts with a step size 1 and uses a reduction factor of 1/2. The CG method is stopped with a relative tolerance of 10^{-6} or after a maximal number of 50 iterations is reached. We initialize with $p^0 = 0$. The homotopy loop starts from $\gamma = 10^{-2}$, reducing γ by a factor of 10 as long as the Newton method converges in a prescribed number of maximal 30 iterations; the minimal allowed γ is 10^{-12} . We denote by $\bar{\gamma}$ the resulting smallest γ for which the Newton method converged. Below, we also report on the number τ_1 of control intervals that exhibit perfect switching, i.e., in which at most one control of N is active. Similarly, we define τ_j for $j > 1$ as the number of time points with $d = j$ in the proximal mapping.

For the spatial discretization, piecewise linear finite elements on an unstructured mesh of 725 triangular elements (maximal diameter 0.1) are chosen. The time discretization is carried out with the cG(1) Petrov–Galerkin method (which corresponds to the Crank–Nicolson method) and 201 equidistant time points. The discrete control is chosen as piecewise constant, i.e., there are 200 degrees of freedom per control component. An implementation of the presented approach in MATLAB can be downloaded from <https://github.com/clason/multiswitchingcontrol>.

Figures 2a, 2b and 3a show the optimal controls for $\alpha = 10^{-1}$ and $N = 3$, $N = 5$ and $N = 7$, respectively (where here and below only active (i.e., nonzero) control components are shown; components that are never active are not listed in the legend, i.e., u_2 , u_4 , u_7 in these cases, respectively). All three configurations terminated with $\bar{\gamma} = 10^{-12}$, and the optimal controls exhibit perfect switching. Note that while all control components are discontinuous,

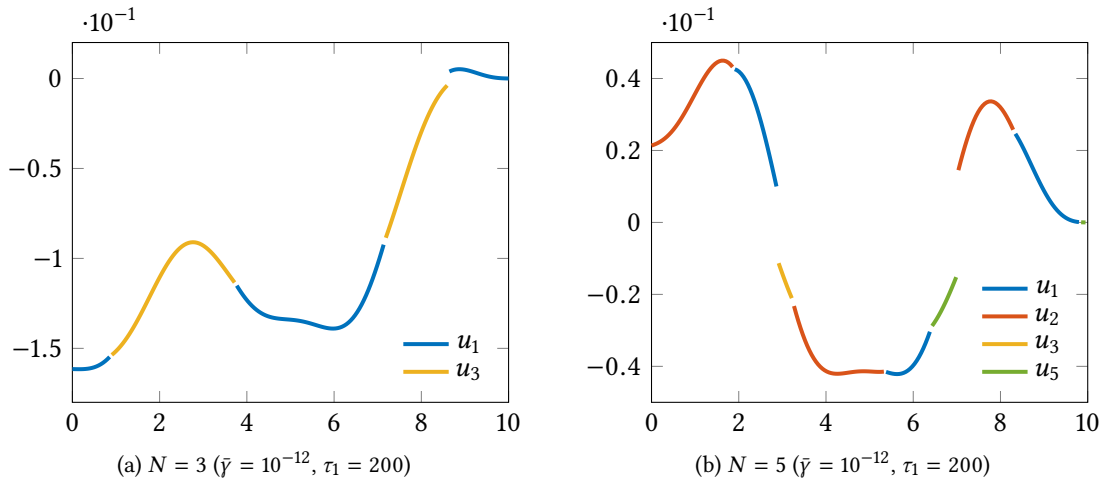


Figure 2: Optimal controls for $\alpha = 10^{-1}$

the norm $\|u(t)\|_{l^1(\mathbb{R}^N)}$ (which corresponds to the absolute value of the curve plotted in the figure) is continuous. This is expected since the L^2 norm of this quantity appears as a penalty in Problem (P).

Controls for $N = 7$ and different values of α are depicted in Figure 3. Since α determines the strength of the switching penalty, it is not surprising that the controls are perfectly switching only for sufficiently large α , as seen in Figure 3a and Figure 3b. For moderate values of α such as $\alpha = 10^{-3}$, there is a single isolated control point where two components are active at the same time (u_3 and u_7 in Figure 3c). For much smaller $\alpha = 10^{-5}$, perfect switching is mostly lost ($\tau_1 = 17$); however, the control still exhibits a switching property, since in most intervals only two and never more than three components are active ($\tau_2 = 140$, $\tau_3 = 43$). Due to the penalization of the L^2 norm in time of the ℓ_1 norm of the controls, we also observe that the magnitude of the plotted control envelope increases with decreasing α while also becoming less regular with respect to time.

The dependence of the switching structure on α is illustrated quantitatively in Table 1, where we also report on the effects on the numerical algorithm. As can be seen, for moderate values of α , the homotopy terminates successfully with $\bar{\gamma} = 10^{-12}$, and both the number of Newton and CG iterations in the final homotopy and Newton step, respectively, increase only slightly. For smaller α and γ , the Newton system becomes increasingly difficult to solve, and the inner CG and Newton iterations terminate due to the maximum number of iterations having been reached, which leads to early termination of the homotopy strategy.

To illustrate the convergence rate of the semismooth Newton method, we report in Table 2 on the convergence history for a single run of the Newton method for $N = 7$, $\alpha = 0.01$ and $\gamma = 10^{-7}$ without homotopy. Both the norm of the residual in the reduced first order optimality condition (4.3) and the number of control intervals where the active control components change (denoted by F_k and s_k , respectively) show superlinear decay. It can also be observed that the reduced optimality conditions are satisfied close to machine precision as soon as the active

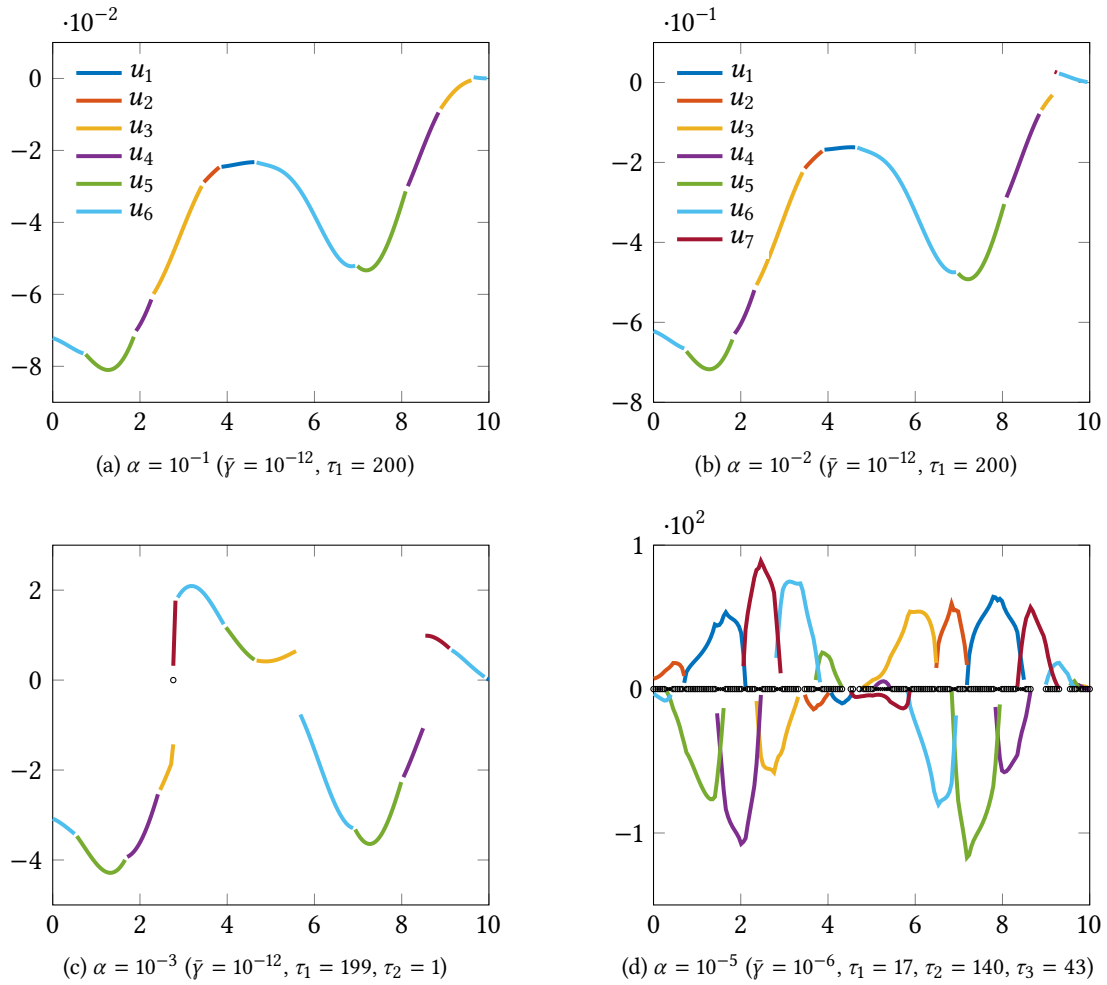


Figure 3: Dependence of the controls on α for $N = 7$. Control intervals without perfect switching are marked on the t -axis

components no longer change. This finite termination of semismooth Newton methods is typical for problems that are linear apart from the case distinction in H_γ and shows the semismooth Newton method's relation to active set methods. The described behavior is typical for large α or γ where no line search is necessary. For moderate values, the iteration first requires some reduced steps before switching to full steps with superlinear convergence and finite termination. For very small α and γ , the prescribed maximal number of iterations is insufficient to reach this convergence region, and the method terminates.

The behavior of the homotopy strategy for moderate $\alpha = 10^{-3}$ is illustrated in Table 3, where both the number of control intervals τ_1, τ_2, τ_3 with one, two or three active control components, respectively, and the number of Newton iterations together with the number of CG iterations in the last Newton step are shown. For large γ , the regularization naturally prevents strong switching, which starts to appear for $\gamma < 10^{-4}$; for $\gamma < 10^{-8}$, the structure of active components

no longer changes. It can also be observed that the number of Newton and CG steps stays relatively constant throughout the whole iteration. The situation is different for small α ; see Table 4. Since the regularized subdifferential depends on the relation between α and γ , the switching structure starts to appear later. Depending on the problem, the limit case for small α, γ might have a solution that is not perfectly switching, in which case the unregularized subdifferential is set-valued. In this example, we observe that for small α and γ , the Newton steps indeed become increasingly difficult to solve, as shown by the growing number of steps necessary before the iteration terminates.

6 Conclusion

Optimal controls for parabolic differential equations with switching structure can be obtained by minimizing a convex functional without needing to introduce additional decision variables or explicit switching points. A combination of Moreau–Yosida regularization and a semismooth Newton method allows the efficient numerical computation of controls for an arbitrary number of control components. Since the approach is based on a general convex analysis framework, it can be adapted to hyperbolic and nonlinear state equations as well as to include sparsity or control constraints.

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Table 1: Dependence on α for $N = 7$ ($\bar{\gamma}$ is the last successful regularization parameter, #SSN is the number of semismooth Newton steps and #CG is the number of CG steps in the last Newton step for $\bar{\gamma}$)

α	10^{-1}	10^{-2}	10^{-3}	10^{-4}	10^{-5}	10^{-6}
τ_1	200	200	199	199	17	2
τ_2	0	0	1	1	140	39
τ_3	0	0	0	0	43	146
$\bar{\gamma}$	10^{-12}	10^{-12}	10^{-12}	10^{-12}	10^{-6}	10^{-7}
#SSN	1	1	4	2	8	15
#CG	3	3	1	1	28	50

Table 2: Convergence history of the semismooth Newton method for $N = 7$, $\alpha = 10^{-2}$ and $\gamma = 10^{-7}$, where F_k denotes the residual norm and s_k denotes the number of control intervals where the active control components change

k	0	1	2	3	4
F_k	$3.133 \cdot 10^{-2}$	$3.131 \cdot 10^{-2}$	$1.902 \cdot 10^{-3}$	$8.285 \cdot 10^{-6}$	$3.463 \cdot 10^{-12}$
s_k	N/A	200	152	4	0

Table 3: Dependence on γ for $N = 7$, $\alpha = 10^{-3}$ (#SSN is the number of semismooth Newton steps, #CG is the number of CG steps in the last Newton step)

γ	10^{-2}	10^{-3}	10^{-4}	10^{-5}	10^{-6}	10^{-8}	10^{-10}	10^{-12}
τ_1	0	0	101	189	198	199	199	199
τ_2	0	45	91	11	2	1	1	1
τ_3	0	101	8	0	0	0	0	0
#SSN	3	4	3	3	3	4	2	4
#CG	4	5	5	4	5	4	1	1

Table 4: Dependence on γ for $N = 7$, $\alpha = 5 \cdot 10^{-5}$ (#SSN is the number of semismooth Newton steps, #CG is the number of CG steps in the last Newton step)

γ	10^{-2}	10^{-3}	10^{-4}	10^{-5}	10^{-6}	10^{-7}	10^{-8}
τ_1	0	0	0	49	169	186	188
τ_2	0	0	0	126	28	14	12
τ_3	0	0	70	19	3	0	0
#SSN	3	3	3	3	7	20	21
#CG	4	6	8	10	11	15	24

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