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Path-following primal-dual interior-point methods for shape optimization of stationary flow problems

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Abstract — We consider shape optimization of Stokes flow in channels where the objective is to design the lateral walls of the channel in such a way that a desired velocity profile is achieved. This amounts to the solution of a PDE constrained optimization problem with the state equation given by the Stokes system and the design variables being the control points of a Bézier curve representation of the lateral walls subject to bilateral constraints. Using a finite element discretization of the problem by Taylor–Hood elements, the shape optimization problem is solved numerically by a path-following primal-dual interior-point method applied to the parameter dependent nonlinear system representing the optimality conditions. The method is an all-at-once approach featuring an adaptive choice of the continuation parameter, inexact Newton solves by means of right-transforming iterations, and a monotonicity test for convergence monitoring. The performance of the adaptive continuation process is illustrated by several numerical examples.

Keywords: PDE constrained optimization, shape optimization, Stokes flow, primal-dual interior-point methods, central path, continuation methods

1. Introduction

Structural optimization problems with constraints given by the Stokes equations can be written as follows

minimize
$$J(\mathbf{u}, p, \alpha)$$
 (1.1a)

over
$$(\mathbf{u}, p, \alpha) \in \mathbf{V} \times Q \times K$$

subject to:
$$S(\mathbf{u}, p, \alpha) = \mathbf{g}$$
. (1.1b)

Here, $J : \mathbf{V} \times Q \times U \rightarrow \mathbb{R}$ denotes the objective functional which depends on the state variables **u** and *p* (the velocity and the pressure) in the state space $\mathbf{V} \times Q$ and

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the design variables α in the design space U. The equation (1.1b) corresponds to the Stokes equations, whereas $K \subset U$ refers to the set of admissible design variables. We note that simplified problems in structural optimization have already been addressed by Bernoulli, Euler, Lagrange and Saint-Venant. However, it became its own discipline during the second half of the last century when the rapidly growing performance of computing platforms and the simultaneously achieved significant improvement of algorithmic tools enabled the appropriate treatment of complex problems (cf. [1-3,13,14,24,25,28,33,35,36,39] and the references therein). The design criteria in structural optimization are determined by a goal oriented operational behavior of the devices and systems under consideration and typically occur as a nonlinear, often non convex, objective functionals which depend on the state variables describing the operational mode and the design variables determining the shape and the topology. The state variables have to satisfy differential equations or systems thereof representing the underlying physical laws. Technological aspects are taken into account by constraints on the state and/or design variables which may occur both as equality and inequality constraints in the model.

The discretization of such structural optimization problems typically gives rise to equality and inequality constrained nonlinear programming problems. If Newton's method is applied to the KKT conditions, each Newton step requires the solution of a linear algebraic system representing the optimality conditions of a related quadratic programming (QP) problem. Hence, Newton methods can be interpreted in the framework of sequential quadratic programming (SOP) which is the most successful method for solving constrained nonlinear optimization problems [4–9,21,22,34]. As far as the appropriate treatment of the inequality constraints is concerned, a local optimum can be approximated from within the feasible set, which is the idea behind interior-point methods. The so-called interior-point revolution in continuous optimization started in the eighties of the last century with Karmarkar's polynomial-time linear programming algorithm. It was immediately found that there is a close relationship to barrier functions which had been used long time before for inequality constrained nonlinear programming problems. Nowadays, interior-point methods are well established tools for constrained nonlinear optimization problems (cf., e.g., [12,15,17-19,20,26,27,29-32,40,41,43-45]). In terms of the barrier parameter, we are faced with a parameter dependent nonlinear system whose solution is referred to as the central path. For the parameter-dependent nonlinear system, continuation methods along the central path are the methods of choice. Such continuation methods have been studied extensively in the literature. A good reference both for theoretical and algorithmic aspects is the recent textbook [16]. The most widely used techniques are path-following predictor-corrector strategies.

In this paper, we consider path-following primal-dual interior-point methods for the shape optimization of stationary flow problems as described by the Stokes system. In particular, we consider Stokes flow in channels where the objective is to design the lateral walls in such a way that a desired velocity profile is obtained. The design variables are chosen as the control points of a Bézier curve representation of the lateral walls. The discretization of the problem is done by Taylor–Hood



Figure 1. Stokes flow in a channel with a backward facing step.

elements with respect to a shape regular family of simplicial triangulations of the computational domain. The resulting optimality conditions for the discretized problem are solved by an adaptive continuation method in the spirit of [16] applied to the parameter dependent nonlinear system representing the central path. For the computation of the derivatives occurring in the KKT conditions and the Hessian, we use automatic differentiation [23].

The paper is organized as follows: In Section 2, we formulate the shape optimization problem along with its finite element discretization and set up the optimality conditions. Section 3 is devoted to the primal-dual approach featuring parameterized minimization subproblems in terms of logarithmic barrier functions that take care of the bilateral constraints on the design variables. The optimality conditions lead to a parameter dependent nonlinear system whose solution is known as the central path. This system is solved by a path-following predictor–corrector type continuation method with an adaptive choice of the continuation parameter (barrier parameter) and a monotonicity test for convergence monitoring. At each continuation step, the Newton increments are computed by means of an iterative scheme based on right-transforming iterations whose construction essentially takes advantage of the specific structure of the associated Hessian. The final Section 4 contains a detailed documentation of numerical examples illustrating the performance of the adaptive continuation method.

2. The shape optimization problem

2.1. Setting of the problem

We consider Stokes flow in a channel occupying bounded domain $\Omega(\alpha) \subset \mathbb{R}^2$ with boundary $\Gamma(\alpha) = \Gamma_{in}(\alpha) \cup \Gamma_{lat}(\alpha) \cup \Gamma_{out}(\alpha)$, where $\Gamma_{lat}(\alpha) := \Gamma_{bot}(\alpha) \cup \Gamma_{top}(\alpha)$, depending on a design variable $\alpha = (\alpha_1, \dots, \alpha_m)^T \in \mathbb{R}^m$ (cf. Fig. 1).

Denoting the velocity and the pressure by $\mathbf{u} := (u_1, u_2)^T$ and p, respectively, and given a velocity and pressure profiles \mathbf{u}^d , p^d , the objective is to design the geometry of the domain in such a way that \mathbf{u}^d , p^d are attained as closely as possible. This

amounts to the minimization problem

$$\inf_{\mathbf{u},p,\alpha} J(\mathbf{u},p,\alpha), \quad J(\mathbf{u},p,\alpha) := \frac{\lambda_1}{2} \int_{\Omega(\alpha)} |\mathbf{u} - \mathbf{u}^d|^2 \, \mathrm{d}x + \frac{\lambda_2}{2} \int_{\Omega(\alpha)} |p - p^d|^2 \, \mathrm{d}x \quad (2.1)$$

subject to the PDE constraint (Stokes flow)

$$-\nu\Delta \mathbf{u} + \nabla p = \mathbf{f} \quad \text{in } \Omega(\alpha) \tag{2.2a}$$

$$\nabla \cdot \mathbf{u} = 0 \quad \text{in } \Omega(\alpha) \tag{2.2b}$$

$$\mathbf{n} \cdot \mathbf{u} = \begin{cases} u_{\text{in}} & \text{on } \Gamma_{\text{in}}(\alpha) \\ u_{\text{out}} & \text{on } \Gamma_{\text{out}}(\alpha) \\ 0 & \text{on } \Gamma_{\text{lat}}(\alpha) \end{cases}$$
(2.2c)

$$\mathbf{t} \cdot \mathbf{u} = 0 \quad \text{on } \Gamma(\alpha) \tag{2.2d}$$

and subject to the bilateral constraints on the design variable α

$$\alpha \in K := \{ \alpha_i \mid \alpha_i^{\min} \leqslant \alpha_i \leqslant \alpha_i^{\max}, \ i = 1, \dots, m \}$$
(2.3)

where $\lambda_i \ge 0$, i = 1, 2, $\lambda_1 + \lambda_2 > 0$, are weighting factors, v is the viscosity, **f** refers to a forcing term, **n**, **t** denote the unit outward normal and unit tangential vector on $\Gamma(\alpha)$, u_{in} , u_{out} stand for a inflow/outflow profiles satisfying the compatibility condition $\int_{\Gamma_{\text{in}}(\alpha)} u_{\text{in}} ds = \int_{\Gamma_{\text{out}}(\alpha)} u_{\text{out}} ds$, and α^{\min} , $\alpha^{\max} \in \mathbb{R}^m$. Note that $u_{\text{out}} := \mathbf{n} \cdot \mathbf{u}$ on $\Gamma_{\text{out}}(\alpha)$ is then determined by means of (2.2b). The state equations (2.2a)–(2.2d) have to be understood in a weak sense. We adopt standard notation from Lebesgue and Sobolev space theory. In particular, we set

$$\mathbf{V} := \{ \mathbf{v} \in H^1(\Omega(\alpha))^2 \mid \mathbf{v} \text{ satisfies (2.2c), (2.2d)} \}, \quad Q := L^2_0(\Omega(\alpha)) .$$

We denote by \mathbf{V}^* and Q^* the duals of \mathbf{V} and Q, and we introduce $A(\alpha) : \mathbf{V} \to \mathbf{V}^*$ and $B(\alpha) : \mathbf{V} \to Q^*$ as the operators associated with the bilinear forms

$$a(\mathbf{v},\mathbf{w};\alpha)) := \mathbf{v} \int_{\Omega(\alpha)} \nabla \mathbf{v} \cdot \nabla \mathbf{w} \, \mathrm{d}x, \quad \mathbf{v} \in \mathbf{V}, \quad \mathbf{w} \in H^1_0(\Omega(\alpha))^2$$
(2.4a)

$$b(\mathbf{v},q;\alpha) := -\int_{\Omega(\alpha)} \nabla \cdot \mathbf{v}q \, \mathrm{d}x, \quad \mathbf{v} \in \mathbf{V}, \quad q \in L^2(\Omega(\alpha)) \,. \tag{2.4b}$$

We further introduce the state space $\mathbf{Y} := \mathbf{V} \times Q$ and the operator $S : \mathbf{Y} \times K \to \mathbf{Y}^*$ according to

$$S(\mathbf{y},\alpha) := \begin{pmatrix} A(\alpha) & B^*(\alpha) \\ B(\alpha) & 0 \end{pmatrix} \begin{pmatrix} \mathbf{u} \\ p \end{pmatrix}, \quad \mathbf{y} = (\mathbf{u},p)^T \in Y, \quad \alpha \in K.$$
(2.5)

We note that $S(\cdot, \alpha) : \mathbf{Y} \to \mathbf{Y}^*, \alpha \in K$, is the Stokes operator. Then, the state equations (2.2a)–(2.2d) can be written in operator form according to

$$S(\mathbf{y},\alpha) = \mathbf{g} \tag{2.6}$$

where $\mathbf{g} := (\mathbf{f}, 0)^T$. The design variables α_i , i = 1, ..., m, are chosen as the Bézier control points in a Bézier representation of $\Gamma(\alpha)$ as a globally continuous composition of polynomial curve segments of some polynomial degree. We choose $\hat{\alpha} \in K$ as a reference design and refer to $\hat{\Omega} := \Omega(\hat{\alpha})$ as the associated reference domain. Then, the actual domain $\Omega(\alpha)$ can be obtained from the reference domain $\hat{\Omega}$ by means of an isomorphism

$$\Omega(\alpha) = \Phi(\hat{\Omega}; \alpha)$$

$$\Phi(\hat{x}; \alpha) = (\Phi_1(\hat{x}; \alpha), \Phi_2(\hat{x}; \alpha))^T, \quad \hat{x} = (\hat{x}_1, \hat{x}_2)^T$$
(2.7)

with continuous components Φ_i , i = 1, 2. We assume: The functions Φ_i , i = 1, 2, are subdifferentiable in $\hat{x} \in \hat{\Omega}$ with uniformly bounded subdifferentials $\partial \Phi_i$ and twice continuously differentiable in α_i , i = 1, ..., m.

The advantage of using the reference domain $\hat{\Omega}$ is that finite element approximations of (2.1) can be performed with respect to that fixed domain without being forced to remesh.

We denote by $(\mathscr{T}_h(\hat{\Omega}))_{\mathbb{N}}$ a shape regular family of simplicial triangulations of $\hat{\Omega}$. By means of (2.7), these triangulations induce an associated family $(\mathscr{T}_h(\Omega(\alpha)))_{\mathbb{N}}$ of simplicial triangulations of the actual physical domains $\Omega(\alpha)$.

We use Taylor–Hood P2/P1 elements (cf., e.g., [11]) for the discretization of the velocity $\mathbf{u} \in \mathbf{V}$ and the pressure $p \in Q$ denoting the associated trial spaces by \mathbf{V}_h and Q_h with dim $\mathbf{V}_h = n_1$ and dim $Q_h = n_2$, respectively. This gives rise to an objective functional $J_h : \mathbb{R}^n \times \mathbb{R}^m$, $n := n_1 + n_2$, by means of

$$J_h(\mathbf{u}_h, p_h, \alpha) := \frac{\lambda_1}{2} (\mathbf{u}_h - \mathbf{u}_h^d)^T I_{1,h}(\alpha) (\mathbf{u}_h - \mathbf{u}_h^d) + \frac{\lambda_2}{2} p_h^T I_{2,h}(\alpha) p_h \qquad (2.8)$$

where $I_{\mathbf{v},h}(\alpha)$, $\mathbf{v} = 1, 2$, are the associated mass matrices and $\mathbf{u}_h^d \in \mathbb{R}^{n_1}$, $p_h \in \mathbb{R}^{n_2}$ result from the L^2 -projections of \mathbf{u}^d , p^d onto \mathbf{V}_h and Q_h , respectively. The discretized shape optimization problem can be stated as

$$\inf_{\mathbf{u}_h, p_h, \alpha} J_h(\mathbf{u}_h, p_h, \alpha) \tag{2.9}$$

subject to the algebraic saddle point problem

$$S_{h}(\mathbf{y}_{h},\alpha) := \begin{pmatrix} A_{h}(\alpha) & B_{h}^{T}(\alpha) \\ B_{h}(\alpha) & 0 \end{pmatrix} \begin{pmatrix} \mathbf{u}_{h} \\ p_{h} \end{pmatrix} = \begin{pmatrix} \mathbf{g}_{1,h} \\ \mathbf{g}_{2,h} \end{pmatrix}$$
(2.10)

representing the Taylor–Hood approximation of the Stokes system (2.5).

Due to the dependence of the domain on the design parameters α_i , i = 1, ..., m, the objective functional J_h is nonconvex. Therefore, there may exist a multitude of local minima. Throughout the following, we assume that $(\mathbf{y}_h^*, \alpha^*) \in \mathbb{R}^n \times K$ is a strict local solution of (2.9), i.e., there exists a neighborhood $\mathscr{U}(\mathbf{y}_h^*, \alpha^*) \subset \mathbb{R}^n \times K$ such that

$$J_h(\mathbf{y}_h^*, \alpha^*) < J(\mathbf{y}_h, \alpha), \quad (\mathbf{y}_h, \alpha) \in \mathscr{U}(\mathbf{y}_h^*, \alpha^*) \setminus \{(\mathbf{y}_h^*, \alpha^*)\}.$$
(2.11)

For notational convenience, in the sequel we will drop the discretization index h.

2.2. Optimality conditions

We introduce Lagrange multipliers $\boldsymbol{\lambda} := (\boldsymbol{\lambda}_{\mathbf{u}}, \boldsymbol{\lambda}_p)^T \in \mathbb{R}^{n_1} \times \mathbb{R}^{n_2}$ and $\boldsymbol{\sigma} := (\sigma_1, \sigma_2)^T$, $\sigma_k \in \mathbb{R}^m_+, k = 1, 2$. The saddle point formulation of the minimization problem (2.9) is

$$\inf_{\mathbf{y},\alpha} \sup_{\boldsymbol{\lambda},\sigma} L(\mathbf{y},\alpha,\boldsymbol{\lambda},\sigma) .$$
 (2.12)

Here, the Lagrangian L is given by

$$L(\mathbf{y}, \alpha, \boldsymbol{\lambda}, \sigma) := J(\mathbf{u}, p, \alpha) + \boldsymbol{\lambda}^{T} (S(\mathbf{y}, \alpha) - \mathbf{g}) + \sigma_{1}^{T} (\alpha^{\min} - \alpha) + \sigma_{2}^{T} (\alpha - \alpha^{\max}).$$
(2.13)

Denoting by $\mathbf{x} := (\mathbf{y}, \alpha)^T$ the primal variables, the first order necessary optimality conditions are given by

$$L_{\mathbf{y}}(\mathbf{x},\boldsymbol{\lambda},\boldsymbol{\sigma}) = 0 \tag{2.14a}$$

$$L_{\alpha}(\mathbf{x}, \boldsymbol{\lambda}, \boldsymbol{\sigma}) = 0 \tag{2.14b}$$

$$L_{\mathbf{y}}(\mathbf{x}, \boldsymbol{\lambda}, \boldsymbol{\sigma}) = 0 \qquad (2.14a)$$

$$L_{\alpha}(\mathbf{x}, \boldsymbol{\lambda}, \boldsymbol{\sigma}) = 0 \qquad (2.14b)$$

$$L_{\boldsymbol{\lambda}}(\mathbf{x}, \boldsymbol{\lambda}, \boldsymbol{\sigma}) = 0 \qquad (2.14c)$$

$$L_{\sigma_{\mathrm{I}}}(\mathbf{x}, \boldsymbol{\lambda}, \boldsymbol{\sigma}) = \alpha^{\mathrm{min}} - \alpha \leq 0 \qquad (2.14d)$$

$$L_{\sigma_{1}}(\mathbf{x}, \boldsymbol{\lambda}, \boldsymbol{\sigma}) = \alpha^{\min} - \alpha \leqslant 0 \qquad (2.14d)$$
$$(\sigma_{1})^{T} (\alpha^{\min} - \alpha) = 0$$

$$L_{\sigma_2}(\mathbf{x}, \boldsymbol{\lambda}, \boldsymbol{\sigma}) = \alpha - \alpha^{\max} \leqslant 0$$

$$(\sigma_2)^T (\alpha - \alpha^{\max}) = 0.$$
(2.14e)

In particular, we find

$$L_{\mathbf{u}}(\cdots) = \lambda_1 I_1(\alpha) (\mathbf{u} - \mathbf{u}^d) + A(\alpha) \lambda_{\mathbf{u}} + B^T(\alpha) \lambda_p = 0$$
(2.15a)

$$L_p(\dots) = \lambda_2 I_2(\alpha) p + B(\alpha) \boldsymbol{\lambda}_{\mathbf{u}} = 0$$
(2.15b)

$$L_{\alpha}(\dots) = \nabla_{\alpha} J(\mathbf{u}, p, \alpha) + \nabla_{\alpha} \langle A(\alpha) \mathbf{u} + B^{T}(\alpha) p - \mathbf{f}, \boldsymbol{\lambda}_{\mathbf{u}} \rangle$$

$$+ \nabla_{\alpha} \langle B(\alpha) \mathbf{u}, \boldsymbol{\lambda}_{p} \rangle - \sigma_{1} + \sigma_{2} = 0$$
(2.15c)

$$L_{\lambda_{\mathbf{u}}}(\cdots) = A(\alpha)\mathbf{u} + B^{T}(\alpha)p - \mathbf{f} = 0$$
(2.15d)

$$L_{\lambda_p}(\cdots) = B(\alpha)\mathbf{u} = 0.$$
 (2.15e)

Theorem 2.1. Let $(\mathbf{y}^*, \alpha^*) \in \mathbb{R}^n \times K$ be a strict local solution of (2.9) and assume that the constraints are qualified in the sense of the linear independence constraint qualification (cf., e.g, [10]). Then, there exist multipliers $\mathbf{\lambda}^* = (\mathbf{\lambda}^*_{\mathbf{u}}, \mathbf{\lambda}^*_p) \in \mathbb{R}^n$ and $\sigma^* = (\sigma^*_1, \sigma^*_2)$, $\sigma^*_k \in \mathbb{R}^m_+$, k = 1, 2, such that $(\mathbf{y}^*, \alpha^*, \mathbf{\lambda}^*, \sigma^*)$ satisfies the KKT system (2.14a)–(2.14e).

Proof. The proof follows from standard results of optimization theory (cf., e.g., [10]).

For the second order derivatives of the Lagrangian L with respect to the primal variables **y** we obtain

$$L_{\mathbf{u}\mathbf{u}}(\cdots) = \lambda_1 I_1(\alpha), \qquad \qquad L_{\mathbf{u}p}(\cdots) = 0 \qquad (2.16a)$$

$$L_{p\mathbf{u}}(\cdots) = 0, \qquad \qquad L_{pp}(\cdots) = \lambda_2 I_2(\alpha). \qquad (2.16b)$$

Consequently, the Hessian L_{xx} has the form

$$L_{\mathbf{xx}}(\cdots) = \begin{pmatrix} I(\alpha) & L_{\mathbf{y}\alpha} \\ L_{\alpha\mathbf{y}} & L_{\alpha\alpha} \end{pmatrix}$$
(2.17)

where

$$I(\alpha) = \begin{pmatrix} \lambda_1 I_1(\alpha) & 0\\ 0 & \lambda_2 I_2(\alpha) \end{pmatrix}$$
(2.18a)

$$L_{\alpha \mathbf{y}} = (L_{\alpha \mathbf{u}} L_{\alpha p}), \qquad L_{\mathbf{y}\alpha} = (L_{\mathbf{u}\alpha} L_{p\alpha})^T.$$
 (2.18b)

Note that

$$L_{\mathbf{u}\alpha} : \mathbb{R}^m \to \mathbb{R}^{n_1}, \quad L_{\alpha \mathbf{u}} : \mathbb{R}^{n_1} \to \mathbb{R}^m$$
(2.19a)

$$L_{p\alpha} : \mathbb{R}^m \to \mathbb{R}^{n_2}, \quad L_{\alpha p} : \mathbb{R}^{n_2} \to \mathbb{R}^m$$
 (2.19b)

$$L_{\mathbf{u}\alpha} = L_{\alpha\mathbf{u}}^*, \quad L_{p\alpha} = L_{\alpha p}^* \tag{2.19c}$$

so that $L_{xx}(\cdots)$ is symmetric. The second order sufficient optimality requires positive definiteness of $L_{xx}(\cdots)$ at optimality.

Theorem 2.2. Let $(\mathbf{x}^*, \boldsymbol{\lambda}^*, \sigma^*) \in \mathbb{R}^n \times K \times \mathbb{R}^n \times (\mathbb{R}^m_+)^2$ satisfy the KKT conditions (2.14a)–(2.14e) and suppose that $L_{\mathbf{xx}}(\mathbf{x}^*, \boldsymbol{\lambda}^*, \sigma^*)$ is positive definite, i.e.,

$$\mathbf{w}^{T}L_{\mathbf{x}\mathbf{x}}(\mathbf{x}^{*},\boldsymbol{\lambda}^{*},\sigma^{*})\mathbf{w} > 0, \quad \mathbf{w} \in \mathbb{R}^{n} \times K \setminus \{\mathbf{0}\}.$$
(2.20)

Then, $(\mathbf{x}^*, \boldsymbol{\lambda}^*, \sigma^*)$ *is a strict local solution of* (2.1).

Proof. We refer to [10].

3. The primal-dual interior-point approach

3.1. The central path

We couple the inequality constraints (2.3) by logarithmic barrier functions with a barrier parameter $\beta = 1/\alpha > 0$, $\alpha \to \infty$, resulting in the following parameterized family of minimization subproblems

$$\inf_{\mathbf{y},\alpha} B^{(\alpha)}(\mathbf{y},\alpha) \tag{3.1}$$

subject to (2.6), where

$$B^{(\alpha)}(\mathbf{y},\alpha) := J(\mathbf{y},\alpha) - \frac{1}{\alpha} \sum_{i=1}^{m} [\ln(\alpha_i - \alpha_i^{\min}) + \ln(\alpha_i^{\max} - \alpha_i)].$$
(3.2)

The dual aspect is to couple the PDE constraint (2.6) by a Lagrange multiplier $\boldsymbol{\lambda} = (\boldsymbol{\lambda}_{\mathbf{u}}, \lambda_p)^T \in \mathbf{V} \times Q$ which gives rise to the saddle point problem

$$\inf_{\mathbf{y},\alpha} \sup_{\boldsymbol{\lambda}} L^{(\alpha)}(\mathbf{y},\boldsymbol{\lambda},\alpha)$$
(3.3)

where the Lagrangian $L^{(\alpha)}$ is given by

$$L^{(\alpha)}(\mathbf{y},\boldsymbol{\lambda},\alpha) = B^{(\alpha)}(\mathbf{y},\alpha) + \langle S(\mathbf{y},\alpha) - \mathbf{g},\boldsymbol{\lambda} \rangle .$$
(3.4)

The central path $\alpha \mapsto x(\alpha) := (\mathbf{y}(\alpha), \boldsymbol{\lambda}(\alpha), \alpha(\alpha))^T$ is given as the solution of the nonlinear system

$$F(x(\alpha), \alpha) = \begin{pmatrix} L_{\mathbf{y}}^{(\alpha)}(\mathbf{y}, \boldsymbol{\lambda}, \alpha) \\ L_{\boldsymbol{\lambda}}^{(\alpha)}(\mathbf{y}, \boldsymbol{\lambda}, \alpha) \\ L_{\alpha}^{(\alpha)}(\mathbf{y}, \boldsymbol{\lambda}, \alpha) \end{pmatrix} = 0$$
(3.5)

which represents the first order necessary optimality conditions for (3.1).

Setting $\mathbf{y}^d := (\mathbf{u}^d, p^d)$, the derivatives of the Lagrangian are given by

$$L_{\mathbf{y}}^{(\alpha)}(\mathbf{y}, \boldsymbol{\lambda}, \alpha) = I(\alpha)(\mathbf{y} - \mathbf{y}^d) + S(\boldsymbol{\lambda}, \alpha)$$
(3.6a)

$$L_{\boldsymbol{\lambda}}^{(\infty)}(\mathbf{y},\boldsymbol{\lambda},\alpha) = S(\mathbf{y},\alpha) - \mathbf{g}$$
(3.6b)

$$L_{\alpha_{i}}^{(\alpha)}(\mathbf{y},\boldsymbol{\lambda},\alpha) = \partial_{\alpha_{i}}J(\mathbf{u},p) + \langle \partial_{\alpha_{i}}S(\mathbf{y},\alpha),\boldsymbol{\lambda} \rangle - \frac{1}{\alpha(\alpha_{i}-\alpha_{i}^{\min})} + \frac{1}{\alpha(\alpha_{i}^{\max}-\alpha_{i})}.$$
(3.6c)

Theorem 3.1. Let $(\mathbf{y}^*, \alpha^*, \lambda^*, \sigma^*) \in \mathbf{Y} \times K \times \mathbb{R}^n \times (\mathbb{R}^m_+)^2$ satisfy the first order necessary optimality conditions (2.14a)–(2.14e) with strict complementarity in (2.14d), (2.14e) and the second order sufficient optimality condition (2.20). Moreover, assume that the linear independence constraint qualification holds true. Then, there exists $\alpha_{\min} > 0$ such that for all $\alpha > \alpha_{\min}$ the minimization subproblems (3.1) admit unique solutions $(\mathbf{y}(\alpha), \boldsymbol{\lambda}(\alpha), \alpha(\alpha))$ satisfying (3.5) and converging to $(\mathbf{y}^*, \boldsymbol{\lambda}^*, \alpha^*)$ as $\alpha \to \infty$.

Proof. We refer to [43].

3.2. Adaptive path-following continuation method

For the solution of the parameter-dependent nonlinear system (3.5) we use an adaptive path-following predictor-corrector strategy along the lines of [16].

Predictor step. The predictor step relies on tangent continuation along the trajectory of the Davidenko equation

$$F_{\mathbf{x}}(\mathbf{x}(\alpha), \alpha) \mathbf{x}'(\alpha) = -F_{\alpha}(\mathbf{x}(\alpha), \alpha) .$$
(3.7)

Given some approximation $\tilde{\mathbf{x}}(\alpha_k)$ at $\alpha_k > 0$, compute $\tilde{\mathbf{x}}^{(0)}(\alpha_{k+1})$, where $\alpha_{k+1} = \alpha_k + \Delta \alpha_k^{(0)}$, according to

$$F_{x}(\tilde{\mathbf{x}}(\alpha_{k}), \alpha_{k}) \ \delta \mathbf{x}(\alpha_{k}) = -F_{\alpha}(\tilde{\mathbf{x}}(\alpha_{k}), \alpha_{k})$$
(3.8a)

$$\tilde{\mathbf{x}}^{(0)}(\boldsymbol{\alpha}_{k+1}) = \tilde{\mathbf{x}}(\boldsymbol{\alpha}_{k}) + \Delta \boldsymbol{\alpha}_{k}^{(0)} \,\delta \mathbf{x}(\boldsymbol{\alpha}_{k}) \,. \tag{3.8b}$$

We use $\Delta x_0^{(0)} = \Delta x_0$ for some given initial step size Δx_0 , whereas for $k \ge 1$ the predicted step size $\Delta x_k^{(0)}$ is chosen by

$$\Delta \alpha_{k}^{(0)} := \left(\frac{\|\Delta \mathbf{x}^{(0)}(\alpha_{k})\|}{\|\tilde{\mathbf{x}}(\alpha_{k}) - \tilde{\mathbf{x}}^{(0)}(\alpha_{k})\|} \frac{\sqrt{2} - 1}{2\Theta(\alpha_{k})} \right)^{1/2} \Delta \alpha_{k-1}$$
(3.9)

where $\Delta \alpha_{k-1}$ is the computed continuation step size, $\Delta \mathbf{x}^{(0)}(\alpha_k)$ is the first Newton correction (see below), and $\Theta(\alpha_k) < 1$ is the contraction factor associated with a successful previous continuation step.

Corrector step. As a corrector, we use Newton's method applied to equation $F(\mathbf{x}(\alpha_{k+1}), \alpha_{k+1}) = 0$ with $\tilde{\mathbf{x}}^{(0)}(\alpha_{k+1})$ from (3.8) as a start vector. In particular, for $\ell \ge 0$ and $j_{\ell} \ge 0$ we compute $\Delta x^{(j_{\ell})}(\alpha_{k+1})$ according to

$$F'(\tilde{x}^{(j_{\ell})}(\alpha_{k+1}), \alpha_{k+1}) \Delta x^{(j_{\ell})}(\alpha_{k+1}) = -F(\tilde{x}^{(j_{\ell})}(\alpha_{k+1}), \alpha_{k+1})$$
(3.10)

and $\overline{\Delta x}^{(j_{\ell})}(\infty_{k+1})$ as the associated simplified Newton correction

$$F'(\tilde{x}^{(j_{\ell})}(\alpha_{k+1}), \alpha_{k+1})\overline{\Delta x}^{(j_{\ell})}(\alpha_{k+1}) = -F(\tilde{x}^{(j_{\ell})}(\alpha_{k+1}) + \Delta x^{(j_{\ell})}(\alpha_{k+1}), \alpha_{k+1}).$$
(3.11)

We monitor convergence of Newton's method by means of

$$\Theta^{(j_{\ell})}(\mathfrak{a}_{k+1}) := \left\| \overline{\Delta x}^{(j_{\ell})}(\mathfrak{a}_{k+1}) \right\| / \left\| \Delta x^{(j_{\ell})}(\mathfrak{a}_{k+1}) \right\|.$$

In case of successful convergence, we accept the current step size and proceed with the next continuation step. However, if the monotonicity test

$$\Theta^{(j_\ell)}(\mathscr{A}_{k+1}) < 1 \tag{3.12}$$

fails for some $j_{\ell} \ge 0$, the continuation step has to be repeated with the reduced step size

$$\Delta \alpha_k^{(\ell+1)} := \left(\frac{\sqrt{2}-1}{g(\Theta^{(j_\ell)})}\right)^{1/2} \Delta \alpha_k^{(\ell)}, \qquad g(\Theta) := \sqrt{\Theta+1}-1 \tag{3.13}$$

until we either achieve convergence or for some prespecified lower bound $\Delta \alpha_{\min}$ observe

$$\Delta \alpha_k^{(\ell+1)} < \Delta \alpha_{\min}$$

In the latter case, we stop the algorithm and report convergence failure.

3.3. Inexact Newton method by right-transforming iterations

As inner iterations we use an inexact Newton method featuring right-transforming iterations (cf., e.g., [29–31]). To this end, we introduce the slack variables

$$z_i^{(1)} := \frac{1}{\alpha \left(\alpha_i - \alpha_i^{\min}\right)}, \quad z_i^{(2)} := \frac{1}{\alpha \left(\alpha_i^{\max} - \alpha_i\right)}, \quad i = 1, \dots, m.$$
(3.14)

In the slack variable formulation of the KKT system, a Newton step in the increments

$$\Delta x := (\Delta \mathbf{u}, \Delta p, \Delta \lambda_{\mathbf{u}}, \Delta \lambda_{p}, \Delta \alpha, \Delta z^{(1)}, \Delta z^{(2)})^{T}$$

reads as follows (for notational convenience, in the following the upper index (∞) in the Lagrangian will be dropped):

where

$$\mathbf{g} := (g_1, g_2, g_3, g_4, g_5, D_1 Z_1 e - x^{-1} e, D_2 Z_2 e - x^{-1} e)^T$$

with appropriate g_i , $i = 1, \ldots, 5$, $e := (1, \ldots, 1)^T$ and

$$D_1 := \operatorname{diag}(\alpha_i - \alpha_i^{\min}), \quad D_2 := \operatorname{diag}(\alpha_i^{\max} - \alpha_i), \quad Z_{\nu} := \operatorname{diag}(z_i^{(\nu)}), \quad \nu = 1, 2.$$

Since D_{ν} , Z_{ν} , $\nu = 1, 2$, are diagonal matrices, the slack variables can be easily eliminated by condensation (block Gaussian elimination) which leads to the condensed Hessian system

$$\begin{pmatrix} \lambda_{1}I_{1}(\alpha) & 0 & | & A(\alpha) & B^{T}(\alpha) & | & L_{\mathbf{u},\alpha} \\ 0 & \lambda_{2}I_{2}(\alpha) & | & \mathbf{B}(\alpha) & 0 & | & L_{p,\alpha} \\ \hline & & & & & & & & & & & & & & & \\ A(\alpha) & B^{T}(\alpha) & | & 0 & 0 & | & L_{\boldsymbol{\lambda}_{u},\alpha} \\ B(\alpha) & 0 & | & 0 & 0 & | & L_{\boldsymbol{\lambda}_{p},\alpha} \\ \hline & & & & & & & & & & & & & & & \\ L_{\alpha,\mathbf{u}} & & L_{\alpha,p} & | & L_{\alpha,\boldsymbol{\lambda}_{u}} & L_{\alpha,\boldsymbol{\lambda}_{p}} & | & \tilde{L}_{\alpha,\alpha} \end{pmatrix} \begin{pmatrix} \Delta \mathbf{v} \\ \Delta \lambda_{p} \\ \hline & & & & & \\ -- \\ L_{\alpha,\mathbf{u}} & L_{\alpha,p} & | & L_{\alpha,\boldsymbol{\lambda}_{u}} & L_{\alpha,\boldsymbol{\lambda}_{p}} \end{pmatrix} \begin{pmatrix} \Delta \mathbf{v} \\ \Delta \lambda \\ \hline & & & \\ -- \\ \Delta \alpha \end{pmatrix} = \begin{pmatrix} I(\alpha) & S & | & L_{\mathbf{v},\alpha} \\ S & 0 & | & L_{\boldsymbol{\lambda},\alpha} \\ \hline & & & & & \\ -- & -- & | & & & \\ L_{\alpha,\mathbf{v}} & L_{\alpha,\boldsymbol{\lambda}} & | & \tilde{L}_{\alpha,\alpha} \end{pmatrix} \begin{pmatrix} \Delta \mathbf{v} \\ \Delta \lambda \\ \hline & & & \\ -- \\ \Delta \alpha \end{pmatrix} = - \begin{pmatrix} \mathbf{h}_{1} \\ \mathbf{h}_{2} \\ \hline & & & \\ \mathbf{h}_{3} \end{pmatrix}$$
(3.15)

with appropriate h_i , i = 1, 2, 3, $\mathbf{v} := (\mathbf{u}, p)^T$, $\boldsymbol{\lambda} := (\boldsymbol{\lambda}_{\mathbf{u}}, \boldsymbol{\lambda}_p)^T$ and

$$S := \left(egin{array}{cc} A(lpha) & B^T(lpha) \ B(lpha) & 0 \end{array}
ight)$$
 $ilde{L}_{lpha lpha} := L_{lpha lpha} + D_1^{-1}Z_1 + D_2^{-1}Z_2.$

Denoting by \tilde{S}^{-1} an approximate inverse of *S* (approximate Stokes solver), an approximate inverse $\tilde{\mathcal{A}}^{-1}$ of the first diagonal block \mathcal{A} of the condensed Hessian *K* is given by

$$\mathscr{A} = \left(egin{array}{cc} I(lpha) & S \ S & 0 \end{array}
ight) \implies \widetilde{\mathscr{A}}^{-1} = \left(egin{array}{cc} 0 & \widetilde{S}^{-1} \ \widetilde{S}^{-1} & -\widetilde{S}^{-1}I(lpha)\widetilde{S}^{-1} \end{array}
ight).$$

Hence, a right-transform K_R of K can be obtained according to

$$K_{R} := \begin{pmatrix} I & 0 & | & -\tilde{S}^{-1}L_{\lambda,\alpha} \\ 0 & I & | & -\tilde{S}^{-1}L_{\mathbf{v},\alpha} + \tilde{S}^{-1}I(\alpha)\tilde{S}^{-1}L_{\lambda,\alpha} \\ -- & -- & | & --- - - - - - - - \\ 0 & 0 & | & I \end{pmatrix}.$$

The right transform K_R provides a regular splitting of the condensed Hessian

$$KK_{R} := \begin{pmatrix} I(\alpha) & S & | & L_{\mathbf{v},\alpha} \\ S & 0 & | & L_{\mathbf{\lambda},\alpha} \\ -- & -- & | & -- \\ L_{\alpha,\mathbf{v}} & L_{\alpha,\mathbf{\lambda}} & | & \tilde{L}_{\alpha\alpha} \end{pmatrix} \begin{pmatrix} I & 0 & | & -\tilde{S}^{-1}L_{\mathbf{\lambda},\alpha} \\ 0 & I & | & -\tilde{S}^{-1}L_{\mathbf{v},\alpha} + \tilde{S}^{-1}I(\alpha)\tilde{S}^{-1}L_{\mathbf{\lambda},\alpha} \\ -- & -- & | & -- \\ 0 & 0 & | & I \end{pmatrix} \\ = \underbrace{\begin{pmatrix} I(\alpha) & S & | & 0 \\ S & 0 & | & 0 \\ -- & -- & | & -- \\ L_{\alpha,\mathbf{v}} & L_{\alpha,\mathbf{\lambda}} & | & \tilde{L}_{\alpha\alpha} \end{pmatrix}}_{=:M_{1}} - \underbrace{\begin{pmatrix} 0 & 0 & | & (I - S\tilde{S}^{-1})(I(\alpha)\tilde{S}^{-1}L_{\mathbf{\lambda},\alpha} - L_{\mathbf{v},\alpha}) \\ 0 & 0 & | & (S\tilde{S}^{-1} - I)L_{\mathbf{\lambda},\alpha} \\ -- & -- & | & -- \\ 0 & 0 & | & 0 \end{pmatrix}}_{=:M_{2} \approx 0}$$

where

$$\hat{L}_{\alpha\alpha} := \tilde{L}_{\alpha\alpha} + L_{\alpha,\lambda} \tilde{S}^{-1} I(\alpha) \tilde{S}^{-1} L_{\lambda,\alpha} - L_{\alpha,\mathbf{v}} \tilde{S}^{-1} L_{\lambda,\alpha} - L_{\alpha,\lambda} \tilde{S}^{-1} L_{\mathbf{v},\alpha}.$$

Let $\Delta \phi^* := (\Delta \mathbf{v}^*, \Delta \lambda^*, \Delta \alpha^*)^T$ be the solution of $K \Delta \phi^* = d$. Then, there holds

$$K \Delta \phi^* = d \iff \underbrace{K K_R}_{= M_1 - M_2} K_R^{-1} \Delta \phi^* = d.$$

The regular splitting $KK_R = M_1 - M_2$ induces the iterative scheme

$$K_R^{-1} \,\delta\phi^{(\nu+1)} \,=\, K_R^{-1} \,\delta\phi^{(\nu)} \,+\, M_1^{-1} \,(d \,-\, K \,\delta\phi^{(\nu)})$$

which results in the right transforming iteration

$$\delta\phi^{(\nu+1)} = \delta\phi^{(\nu)} + K_R M_1^{-1} (d - K \,\delta\phi^{(\nu)}). \tag{3.16}$$

Let $\delta \phi^{(0)} := (\delta \mathbf{v}^{(0)}, \delta \boldsymbol{\lambda}^{(0)}, \delta \alpha^{(0)})^T$ be given and compute $\delta \phi^{(\ell+1)}, k \ge 0$, by

$$\delta \phi^{(\ell+1)} = \delta \phi^{(\ell)} + K_R M_1^{-1} (d - K \, \delta \phi^{(\ell)}).$$

Step 1. Compute the residual $r := d - K \delta \phi^{(\ell)}$.

Step 2. Compute $\delta \psi = (\delta \psi_1, \delta \psi_2, \delta \psi_3)^T$ as the solution of

$$M_1 \,\delta\psi = \begin{pmatrix} I(\alpha) & S & \mid 0 \\ S & 0 & \mid 0 \\ -- & -- & \mid -- \\ L_{\alpha,\mathbf{v}} & L_{\alpha,\boldsymbol{\lambda}} & \mid \hat{L}_{\alpha\alpha} \end{pmatrix} \begin{pmatrix} \delta\psi_1 \\ \delta\psi_2 \\ -- \\ \delta\psi_3 \end{pmatrix} = \begin{pmatrix} r_1 \\ r_2 \\ -- \\ r_3 \end{pmatrix}.$$

Step 2.1. Compute $(\delta \psi_1, \delta \psi_2)^T$ according to

$$\begin{pmatrix} \delta\psi_1\\ \delta\psi_2 \end{pmatrix} = \begin{pmatrix} I(\alpha) & S\\ S & 0 \end{pmatrix}^{-1} \begin{pmatrix} r_1\\ r_2 \end{pmatrix} \approx \begin{pmatrix} 0 & \tilde{S}^{-1}\\ \tilde{S}^{-1} & -\tilde{S}^{-1}I(\alpha)\tilde{S}^{-1} \end{pmatrix} \begin{pmatrix} r_1\\ r_2 \end{pmatrix}.$$

Step 2.2. Compute $\delta \psi_3$ as the solution of

$$\hat{L}_{\alpha\alpha}\,\delta\psi_3\,=\,r_3\,-\,L_{\alpha,\mathbf{v}}\delta\psi_1\,-\,L_{\alpha,\mathbf{\lambda}}\delta\psi_2\,.$$

As a termination criterion for the inexact Newton solves, we stop the inner iterations, if

$$\boldsymbol{\vartheta}_{\ell} := \frac{1 - \delta_{\ell+1}}{1 + \overline{\delta}_{\ell}} \frac{\|\boldsymbol{K}_{R}^{-1} \,\delta\boldsymbol{\phi}^{(\ell+1)}\|}{\|\boldsymbol{K}_{R}^{-1} \,\delta\boldsymbol{\phi}^{(\ell)}\|} \leqslant \frac{1}{4}$$
(3.17)

where

$$\overline{\delta}_{\ell+1} := rac{\sqrt{arepsilon_\ell}}{\|\delta \phi^{(\ell)}\|}$$

with ε_{ℓ} being a lower bound for $\|\Delta \phi - \delta \phi^{(\ell)}\|$. Note that

$$K_R^{-1} \,\delta\phi^{(\ell)} = \begin{pmatrix} I & 0 & | & \tilde{S}^{-1}L_{\boldsymbol{\lambda},\alpha} \\ 0 & I & | & \tilde{S}^{-1}L_{\mathbf{v},\alpha} - \tilde{S}^{-1}I(\alpha)\tilde{S}^{-1}L_{\boldsymbol{\lambda},\alpha} \\ \hline & & & \\ 0 & 0 & | & & I \end{pmatrix} \begin{pmatrix} \delta \mathbf{v}^{(\ell)} \\ \delta \boldsymbol{\lambda}^{(\ell)} \\ \hline & & \\ \delta d^{(\ell)} \end{pmatrix}.$$

4. Numerical results

In the numerical examples, we consider Stokes flow with viscosity v in channels $\Omega(\alpha) \subset \mathbb{R}^2$ with reference domain $\hat{\Omega} := \Omega(\alpha^0)$ whose boundaries Γ consist of the inflow and outflow boundaries Γ_{in} , Γ_{out} and the lateral boundaries Γ_{top} , Γ_{bot} (cf. Fig. 1). The inflow u_{in} on the inflow boundary Γ_{in} and the desired profiles \mathbf{u}^d , p^d as well as the weighting factors λ_i , i = 1, 2, are given. We assume m_1 Bézier control points α_i , $i = 1, \ldots, m_1$, on $\hat{\Gamma}_{top}$ and $m_2 = m - m_1$ Bézier control points α_i , $i = m_1 + 1, \ldots, m$, on $\hat{\Gamma}_{bot}$ with given lower and upper bounds α_i^{\min} , α_i^{\max} , $i = 1, \ldots, m$. We use a discretization by Taylor–Hood P2/P1 elements [11] with respect to a simplicial triangulation $\mathcal{T}_h(\hat{\Omega})$ of the reference domain $\hat{\Omega}$ with n_1 degrees of freedom (DOF) for each velocity component and n_2 DOF for the pressure. The derivatives occurring in the KKT conditions and the Hessians have been computed by automatic differentiation [23]. We further refer to $k \in \mathbb{N}$ and $\ell_k \in \mathbb{N}$ as the numbers of the actual continuation step and the inexact Newton iteration, respectively. Moreover, α and $\Delta \alpha$ denote the barrier parameter and its increment, Θ stands for the parameter in the monotonicity test (cf. (3.12)), $\|\Delta \alpha\|_2$ and $\|\alpha - \alpha^*\|_{\infty}$ are the

Table 1.			
Example 1. Discretiza	ation data and	l termination	criteria.

$2n_1$	<i>n</i> ₂	m_1	<i>m</i> ₂	∞^0	$\Delta \infty_0$	tolnt	tolmin	<i>h</i> _{max}
4998	646	4	4	100	200	4.0e-03	1.0e-06	0.15

Table 2.

2

Example 1. Convergence history of the continuation method.

k	x	Δ∝	corr.	ℓ_k	$\ \Delta \alpha\ _2$	$\ \alpha - \alpha^*\ _{\infty}$	J	Θ
0	100.0	(200.0)		_	_	4.0e-01	1.8e+00	
1	100.0	163.3	no	1	6.9e-01	1.4e-03	1.1e-05	0.36
				2	6.0e-06	1.4e-03	1.1e-05	0.33
2	263.3	459.8	no	1	1.9e-03	5.0e-04	1.5e-06	0.03
3	723.1				_	4.0e-04	8.8e-07	

Euclidean norm of the Newton increment in the design variables and the maximum norm of the error in the design variables (in case the optimal design α^* is explicitly known), and J is the value of the objective functional at the *k*-th continuation step. Finally, 'tolmin' and 'tolnt' refer to the termination criteria of the continuation process and the inexact Newton solves, respectively, and h_{max} denotes the maximal mesh width.

Example 1. The data of the first example are as follows:

$$\begin{split} \hat{\Omega} &= \Omega(\alpha^0) = (1,5) \times (-1,+1) \text{ with } \alpha^0 = (+1,+1,+1,+1,-1,-1,-1,-1)^T \\ \mathbf{u}^d, p^d \text{ according to } \overline{\alpha} &= (1.00,1.28,1.40,1.24,-1.00,-1.05,-1.20,-1.13)^T \\ \mathbf{v} &= 1.0 , \ u_{\text{in}}(x_1,x_2) = (6(1+x_2)(1-x_2), \ 0) \\ \alpha^{\text{min}} &= (0,0,0,0,-5,-5,-5,-5)^T, \quad \alpha^{\text{max}} = (5,5,5,5,0,0,0,0)^T. \end{split}$$

The optimal solution $(\mathbf{u}^*, p^*, \alpha^*)$ has been chosen as

$$\mathbf{u}^* = \mathbf{u}^d, \quad p^* = p^d, \quad \alpha^* = \overline{\alpha}$$

where $\overline{\alpha}$ is the design associated with \mathbf{u}^d, p^d . Hence, the value J^* of the objective functional at optimality is zero.

The data of the discretization, the initial value α^0 of the inverse of the barrier parameter and its initial increment $\Delta \alpha^0$ as well as the termination criteria tolnt and tolmin are shown in Table 1. Table 2 contains the convergence history of the adaptive path-following primal-dual interior-point method. The algorithm was stopped by criterion tolmin after three continuation steps and a computational time of 55 minutes (AMD Athlon, 1.8 GHz) with the final design vector

$$\alpha^{3} = (1.0000, 1.2797, 1.4001, 1.2399, -1.0000, -1.0496, -1.2000, -1.1297)^{T}$$

m



Figure 2. Results for Example 1. Initial shape (top left), optimal shape after the third iteration (top right), the optimal velocity field as an arrow plot (bottom left) and the L_0^2 -gauged pressure field shown by equipotential lines (bottom right).

Due to our experience, a fast convergence of the algorithm requires initial parameters α^0 and $\Delta \alpha^0$ that are not chosen too small. In particular, for too small $\Delta \alpha^0$ the algorithm starts too 'carefully and does not achieve satisfactory progress within a reasonable number of outer iterations.

Figure 2 displays the initial and optimal shape (top) and the associated velocity and pressure field (bottom).

Example 2. This example has been constructed in such a way that the desired velocity profile is outside the range of the admissible design variables. Consequently, the value J^* of the objective functional at optimality is greater than zero (cf. Table 4). It can be observed that the upper computed Bézier curve tries to 'mimic' the corresponding part of the shape associated with the desired velocity (cf. Fig. 3, bottom right). The data for α^{\min} , α^{\max} , \mathbf{u}_{in} , etc. are the same as in Example 1, except that the desired velocity field corresponds to the following vector of design variables

 $\bar{\alpha} = (2.50, 1.65, 1.15, 0.90, -1.30, -0.85, -1.20, -1.35)^T$.

After reaching the criterion $\alpha > \alpha_{max} = 15000$, the computed optimal design turned

Table 3.

Example 2. Discretization data and termination criteria.

$2n_1$	<i>n</i> ₂	m_1	<i>m</i> ₂	∞^0	$\Delta \propto^0$	tolnt	tolmin	h _{max}
2762	361	4	4	100	300	4.0e-03	1.0e-06	0.20

Table 4.
Example 2. Convergence history of the continuation method.

k	x	Δ∝	corr.	ℓ_k	$\ \Delta \alpha\ _2$	J	Θ
0	100.0	(300.0)		_		2.4e+01	
1	100.0	269.6	no	1	1.8e+00	4.0e+00	0.30
				2	1.2e-02	3.9e+00	0.98
				3	2.3e-01	1.9e+00	0.51
				4	1.7e-01	7.0e-01	0.25
2	369.6	274.3	no	1	3.1e-02	6.8e-01	0.17
				2	5.8e-03	6.7e-01	0.24
3	643.8	512.2	no	1	1.3e-03	6.7e-01	0.06
4	1156.0	1321.6	no	1	7.4e-04	6.7e-01	0.03
5	2477.6	3205.5	no	1	7.2e-04	6.7e-01	0.04
6	5683.1	1584.6	no	1	4.1e-04	6.7e-01	0.85
7	7267.7	9043.7	no	1	2.1e-04	6.7e-01	0.01
8	16311.4	(8077.4)	no	1	1.3e-04	6.7e-01	0.54

out to be

$$\alpha^8 = (1.0000, 0.5171, 0.0928, 0.0000, -1.0000, -0.6795, -1.2788, -1.0710)^T$$

Example 3. In this example, we consider a channel with a backward facing step (cf. Fig. 4, top left). Here, the three horizontal segments of the geometry are fixed and only the segment that connects the two lower horizontal segments is variable. In this setting, the function that describes the bottom segment of the geometry is a *composite* Bézier curve consisting of three curves of degree 0, 4 and 0, respectively. Note that the composite curve is only continuous. The data for this geometry are as follows:

$$\begin{split} \hat{\boldsymbol{\Omega}} &= \boldsymbol{\Omega}(\alpha^0) \text{ with } \alpha^0 = (+1.0, -0.5, -0.5, -1.5, -1.5, -1.5)^T \\ \mathbf{u}^d, p^d \text{ according to } \overline{\alpha} &= (+1.0, -0.5, -0.55, -0.6, -0.6, -1.5)^T \\ \boldsymbol{\nu} &= 1.0, \ \boldsymbol{u}_{\text{in}}(x_1, x_2) = (6(1+x_2)(1-x_2), \ 0) \\ \boldsymbol{\alpha}^{\text{min}} &= (0, -5, -5, -5, -5, -5)^T, \quad \boldsymbol{\alpha}^{\text{max}} = (5, 0, 0, 0, 0, 0)^T . \end{split}$$

As in Example 1, the optimal solution $(\mathbf{u}^*, p^*, \alpha^*)$ has been chosen as $\mathbf{u}^* = \mathbf{u}^d$, $p^* = p^d$ and $\alpha^* = \overline{\alpha}$. Table 5 provides information about the discretization and Table 6 displays the convergence history of the adaptive path-following algorithm.



Figure 3. Results for Example 2. Initial shape (left) and the optimal shape after 8 iterations (right). Here, the first Bézier point of each curve (top and bottom) of the initial shape are fixed and different from the corresponding points of the shape corresponding to the desired velocity profile (dotted line).



Figure 4. Results for Example 3. A scheme of the backward step geometry (top left) along with the optimal pressure filed (top right), initial and optimal shape after 7 outer iterations (bottom left and right).

Table 5.					
Example 3.	Discretization	data a	and	termination	criteria.

$2n_1$	<i>n</i> ₂	m_1	<i>m</i> ₂	∞^0	$\Delta \infty^0$	tolnt	tolmin
2856	191	1	5	100	300	4.0e-03	1.0e-06

Table 6.

Example 3. Convergence history of the continuation method.

k	x	Δ∝	corr.	ℓ_k	$\ \Delta \alpha\ _2$	$\ \alpha - \alpha^*\ _{\infty}$	J	Θ
0	100.0	(300.0)				9.0e-01	2.6e+00	
1	100.0	(300.0)	yes	1	1.2e+00	1.7e-01	9.6e-01	0.58
				2	8.8e-01	1.7e-01	1.3e-01	618.42
1	100.0	425.5	no	1	2.1e-01	3.3e-02	4.3e-04	0.11
2	525.5	417.1	no	1	1.2e-01	3.3e-02	2.3e-03	0.41
				2	3.3e-02	2.5e-02	2.3e-03	0.58
				3	1.6e-02	2.4e-02	2.0e-03	0.92
				4	2.0e-02	2.9e-03	1.6e-05	0.43
				5	5.7e-04	3.2e-03	2.5e-05	_
3	942.6	323.5	no	1	2.9e-03	3.6e-03	5.1e-05	0.34
4	1266.1	283.7	no	1	1.4e-03	3.5e-03	4.9e-05	0.27
5	1549.8	593.1	no	1	1.7e-04	2.9e-03	3.3e-05	0.05
6	2142.9	2265.3	no	1	1.3e-04	2.1e-03	1.7e-05	0.01
7	4408.2	_	—		_	2.0e-04	1.9e-07	_

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