Adjoint-based computation of shape sensitivity in a Rijke-tube

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Abstract
Geometrical shape is the most important parameter defining acoustic spectra. While adjoint-based sensitivity analysis of shape alterations is a well established technique in other engineering disciplines such as airfoil design, it is less common in acoustic and especially thermoacoustic contexts. Only recently the thermoacoustic community has begun to deploy such methods for stability assessment. The present paper aims at continuing this effort by discussing the challenges of domain shape deformations to adjoint-based gradient calculations and a parameter-free implementation into a state-of-the-art thermoacoustic Helmholtz solver. Deformations of a simple thermoacoustic system – a Rijke tube – will be studied as an example case to validate the method. Despite this practical focus on thermoacoustics, the results of the paper are fruitful for other acoustic topics involving stability analyses, as the underlying concepts readily apply to any eigenvalue problem linear or nonlinear in its eigenvalue. Indeed, the gradient information calculated from first-order theory is the first building block for efficient algorithms to optimize the shape of acoustic devices.

Keywords: adjoints, thermoacoustics, shape derivatives, eigenvalue problems

1 INTRODUCTION
The most influential parameter to acoustic resonance is shape and this paper discusses how to efficiently compute the influence of shape deformations to thermoacoustic stability based on an adjoint method.

In gas turbine engineering a critical aspect is thermoacoustic stability of the combustor. This is the robustness of the combustor against self-excited pressure fluctuations that arise from a constructive feedback of the fluctuations of the rate of heat release in the flame and the acoustics of the combustion chamber. A linear modeling strategy for this phenomenon is the use of a Helmholtz equation that is equipped with a term modeling the acoustic feedback due to heat release [1]. It reads:

\[ \nabla \cdot (c^2 \hat{\nabla} p) + \omega^2 \hat{p} + F(\omega) \nabla \hat{p}_{\text{ref}} \cdot \hat{n}_{\text{ref}} = 0 \quad \text{in } \Omega. \]

(1)

Here, \( \hat{p} \) is the Fourier transform of the pressure fluctuation, \( c \) the local speed of sound, \( \omega \) the complex eigenfrequency, \( F(\omega) \) is a function modeling the flame response linking the velocity fluctuation in the direction \( \hat{n}_{\text{ref}} \) at the point \( \hat{x}_{\text{ref}} \) to the fluctuation in heat release rate in the flame and accounting for the thermodynamic properties of the flame. \( \Omega \) denotes the considered domain. It is confined by its boundary \( \partial \Omega \) at which impedance conditions are set. Mathematically these are mixed-type boundary conditions

\[ D(\omega) \hat{p} + \nabla \hat{p} \cdot \hat{n} = 0 \quad \text{on } \partial \Omega. \]

(2)

where \( D \) is typically defined from a normalized impedance \( Z \) as \( D = \frac{i \omega}{Z} \) and \( \hat{n} \) denotes the outward-pointing unit normal. The equations (1) and (2) present an eigenvalue problem. In contrast to an acoustic Helmholtz equation, the thermoacoustic Helmholtz equation is highly non-linear in its eigenvalue, and non-self-adjoint due to the flame response term and possibly due to the boundary conditions if they are not energy conserving. The review [2] contains a recent overview on nonlinear eigenvalue problems and their mathematics.

How the eigenvalue \( \omega \) depends on the domain \( \Omega \) is the key question discussed in this paper. In particular, how the eigenvalue changes with infinitesimal changes in the domain shape at first order is discussed. Such a change
is commonly referred to as shape derivative:
\[
\frac{\partial \omega}{\partial \varepsilon} = \lim_{\varepsilon \to 0} \frac{\omega(\Omega(\varepsilon)) - \omega(\Omega(0))}{\varepsilon}.
\] (3)

Here, \( \Omega(\varepsilon) \) describes a family of domains smoothly governed by some parameter \( \varepsilon \). An example of such a distortion is depicted in Fig. 1.

Shape derivatives of eigenvalue problems have a long tradition. A classic example is the buckling of a column problem (e.g. [3]). A general mathematical introduction to shape gradients and shape optimization can be found in [4]. Sensitivity of eigenvalue problems is discussed in [5]. In the recent years, the thermoacoustic community has applied adjoint methods to compute the sensitivity of the eigenvalue with respect to many design parameters [6, 7]. Only little attention is paid to the domain shape. For instance, Aguilar et al. used low-order network models for modeling longitudinal modes, i.e., configurations that are essentially one-dimensional. In these models the length of a tube appears as an explicit modeling parameter, such that shape may be treated like any other modeling parameter [8]. This approach can be extended to three dimensions by introducing a parameterized discretization mesh. The parameters of which will then explicitly appear in the considered equations. This is an idea that is used for instance in airfoil optimization using parameterized airfoil shapes [9]. However, the current study follows an approach which does not require to incorporate the considered deformations on the level of a computational grid; in that sense the approach is independent of a parameterized mesh. The mathematics follows the derivation in [10], but is simplified to be independent of the mode normalizations.

2 THEORY

In general, a linear operator \( \mathcal{L} \) consists of differential operators defined in a domain \( \Omega \), with corresponding boundary conditions valid on \( \partial\Omega \). In this report the notation
\[
\mathcal{L}\hat{\rho} := \begin{cases} \mathcal{L}_\Omega \hat{\rho} \text{ in } \Omega \\ \mathcal{L}_{\partial\Omega} \hat{\rho} \text{ on } \partial\Omega \end{cases}
\] (4)

will be used to distinguish between these two components of \( \mathcal{L} \).

Studying the eigenvalue problem
\[
\mathcal{L}(\omega, \varepsilon)\hat{\rho} = 0
\] (5)

where \( \varepsilon \) is a parameter determining the domain shape, i.e., \( \Omega = \Omega(\varepsilon) \), is straightforward by means of perturbation theory. Introducing the ansätz
\[
\omega = \omega_0 + \omega_1 \varepsilon + \text{h.o.t. and } \hat{\rho} = \hat{\rho}_0 + \hat{\rho}_1 \varepsilon + \text{h.o.t.}
\] (6)

and the Taylor series expansion of \( \mathcal{L} \)
\[
\mathcal{L}(\omega, \varepsilon) = \mathcal{L}(\omega_0, 0) + \partial_\omega \mathcal{L}(\omega, 0) \Delta \omega + \partial_\varepsilon \mathcal{L}(\omega_0, 0) \varepsilon + \text{h.o.t.}
\] (7)
into the eigenvalue problem (5), yields
\[ (L_{0,0} + L_{1,0} \omega_1 + L_{0,1} \epsilon) (\hat{p}_0 + \hat{p}_1 \epsilon) + \text{h.o.t.} = 0. \] (8)

Sorting for like powers of \( \epsilon \) gives the first-order equation
\[ L_{0,0} \hat{p}_1 + L_{1,0} \hat{p}_0 \omega_1 + L_{0,1} \hat{p}_0 = 0. \] (9)

This is the standard first-order equation for the eigenvalue problem. The unknowns \( \omega_1 \) and \( \hat{p}_1 \) are found by applying the inner product with the adjoint baseline solution \( \hat{p}_0 \).

The adjoint operator of \( L \) is commonly defined to satisfy
\[ \langle a | L_\Omega b \rangle = \left( L_\Omega^\dagger a \right) b \] + const. \( \forall a, b, \) (10)

where const. originates from boundary integrals that result from an integration by parts. This constant is called boundary conjunct in [10] and denoted in the following as
\[ \text{const.} := [a | b]_{\Omega}, \] (11)

The boundary conjunct is typically used to specify the boundary conditions \( L_\Omega^\dagger \) for the adjoint equation by demanding that the integration constant is vanishing. Such adjoint boundary conditions allow for a convenient usage of the adjoint operator as subsequent calculations do not need to consider the constant. However, as the discussion below shows, the boundary conjunct is required when the boundary itself is subject to a perturbation, i.e., displacements of the boundary are to be studied using perturbation methods.

2.1 Expansion of the boundary conditions at first order

To see the effect of boundary displacement, we assume that for all points on \( \partial \Omega(0) \) there is a mapping to \( \partial \Omega(\epsilon) \). This mapping is assumed to be one-to-one. Hence, locally we have for each point on \( \partial \Omega(0) \) a trajectory \( \bar{x}_0 = \bar{x}_0(\epsilon) \), such that \( \bar{x}_0(\epsilon) \) will be on \( \partial \Omega(\epsilon) \). These trajectories are not unique. Yet, with this notation\(^1\) we can expand a function value for a point \( \bar{x}_0 \):
\[ \hat{p}(\bar{x}_0(\epsilon)) = \hat{p}(\bar{x}_0(0)) + \hat{p}'(\bar{x}_0(0)) \frac{\partial \bar{x}_0}{\partial \epsilon} |_{\epsilon=0} \epsilon + \text{h.o.t.} \] (12)

Using the expansion of a boundary value (12) at first order for a mixed boundary condition (2), we get the following first-order expansion in terms of \( \epsilon \) and \( \omega \):
\[ 0 = D(\omega_1) \left( \hat{p}(\bar{x}_0(0)) + \hat{p}'(\bar{x}_0(0)) \frac{\partial \bar{x}_0}{\partial \epsilon} \right) + \nabla \left( \hat{p}(\bar{x}_0(0)) + \hat{p}'(\bar{x}_0(0)) \frac{\partial \bar{x}_0}{\partial \epsilon} \right) \cdot \left( \bar{n}(\bar{x}_0(0)) + \bar{n}'(\bar{x}_0(0)) \frac{\partial \bar{x}_0}{\partial \epsilon} \right) \]
\[ + D'(\omega_1) \hat{p}(\bar{x}_0(0)) \Delta \omega. \] (13)

At first order, this amounts to:
\[ 0 = D(\omega_1) \hat{p}(\bar{x}_0(0)) + \nabla \hat{p}(\bar{x}_0(0)) \cdot \bar{n}(\bar{x}_0(0)) \]
\[ + \left( D(\omega_1) \hat{p}'(\bar{x}_0(0)) \frac{\partial \bar{x}_0}{\partial \epsilon} + \nabla \left( \hat{p}'(\bar{x}_0(0)) \frac{\partial \bar{x}_0}{\partial \epsilon} \right) \cdot \bar{n}(\bar{x}_0(0)) + \nabla \hat{p}(\bar{x}_0(0)) \cdot \left[ \bar{n}'(\bar{x}_0(0)) \frac{\partial \bar{x}_0}{\partial \epsilon} \right] \right) \epsilon \]
\[ + D'(\omega_1) \hat{p}(\bar{x}_0(0)) \Delta \omega. \] (14)

\(^1\) Please note the usage of ()' for a derivative, which is considered to be a row vector in cases of spatial arguments.
From (9) we find the boundary condition to the first-order correction \( \hat{p}_1(\vec{x}) \):

\[
0 = \mathcal{D}(\omega_0) \hat{p}_1(\vec{x}_0(0)) \nabla \hat{p}_1(\vec{x}_0(0)) \cdot \hat{n}(\vec{x}_0(0)) + \mathcal{D}'(\omega_0) \hat{p}_0(\vec{x}_0(0)) \omega_0 \\
+ \left( \mathcal{D}(\omega_0) \hat{p}_0(\vec{x}_0(0)) \frac{d\vec{x}_0}{d\varepsilon} + \left( \nabla \left[ \hat{p}_0(\vec{x}_0(0)) \frac{d\vec{x}_0}{d\varepsilon} \right] \cdot \hat{n}(\vec{x}_0(0)) + \nabla \hat{p}_0(\vec{x}_0(0)) \cdot \left[ \hat{n}(\vec{x}(0)) \frac{d\vec{x}}{d\varepsilon} \right] \right) \right) .
\]

(15)

All \( x_0 \) are evaluated at \( \varepsilon = 0 \) which means this is a condition holding on the unperturbed surface:

\[
D(\omega_0) \hat{p}_1 + \nabla \hat{p}_1 \cdot \hat{n} = -r_1 - \omega_1 \mathcal{D}'(\omega_0) \hat{p}_0 \quad \text{on} \ \partial \Omega(0)
\]

(16)

Because the perturbed boundary condition is formulated on the unperturbed surface, the approach presented in this paper operates on the unperturbed domain only. This is the very reason why no parameterized mesh that evolves with the perturbation parameter \( \varepsilon \) needs to be created. The deformation enters the problem statement in terms of \( d\vec{x}/d\varepsilon \) included in \( r_1 \).

2.2 The boundary conjunct of wave equations

For the thermoacoustic Helmholtz equation, the boundary conjunct arises from an integration by parts of the Laplacian. It reads

\[
[a,b]_{V(\cdot;V)} = \oint_{\partial \Omega(0)} c^2 \left( n \nabla b \cdot \vec{n} - \nabla a \cdot \vec{n} b \right) dS .
\]

(17)

If \( a = \hat{p}^\ddagger \) and \( b = \hat{p} \), we find

\[
[\hat{p}^\ddagger | \hat{p}]_{V(\cdot;V)} = \oint_{\partial \Omega(0)} c^2 \left( \vec{n} \nabla \hat{p} \cdot \vec{n} - \nabla \hat{p}^\ddagger \cdot \vec{n} \hat{p} \right) dS = -\oint_{\partial \Omega(0)} c^2 \left( \vec{n} \nabla D(\omega_0) \hat{p} - \mathcal{D}' \hat{p} \right) dS = 0
\]

(18)

Here \( D^\ddagger(\omega_0) \) denotes the function determining the adjoint mixed-type boundary condition. In order to make the boundary conjunct vanish, the adjoint boundary condition needs to be specified by \( D^\ddagger(\omega) = \mathcal{D}(\omega) \).

With this specification, we find a particular formula for the boundary conjunct of \( \hat{p}_0 \) with \( \hat{p}_1 \):

\[
[\hat{p}_0 | \hat{p}_1]_{V(\cdot;V)} = \oint_{\partial \Omega(0)} c^2 \left( \hat{p}_0 \nabla \hat{p}_1 \cdot \vec{n} - \nabla \hat{p}_0 \cdot \vec{n} \hat{p}_1 \right) dS
\]

\[
= \oint_{\partial \Omega(0)} c^2 \left( \hat{p}_0 \nabla \hat{p}_1 \cdot \vec{n} - (\mathcal{D}(\omega_0) \hat{p}_0) \hat{p}_1 \right) = \oint_{\partial \Omega(0)} c^2 \hat{p}_0^\ddagger \left( \nabla \hat{p}_1 \cdot \vec{n} + \mathcal{D}(\omega_0) \hat{p}_0 \right)
\]

\[
= \oint_{\partial \Omega(0)} c^2 \hat{p}_0^\ddagger \left( -r_1 - \omega_1 \mathcal{D}'(\omega_0) \hat{p}_0 \right) dS = -\oint_{\partial \Omega(0)} c^2 \hat{p}_0^\ddagger r_1 dS - \omega_1 \oint_{\partial \Omega(0)} c^2 \hat{p}_0^\ddagger \mathcal{D}'(\omega_0) \hat{p}_0 dS
\]

(19)

\[
= -\left< \hat{p}_0^\ddagger | r_1 \right> - \omega_1 \left< \hat{p}_0^\ddagger | \mathcal{D}'(\omega_0) \hat{p}_0 \right> .
\]

Note, that the boundary conjunct consists of surface integrals of quantities that are derived from the known baseline solution \( \hat{p}_0 \) and \( \hat{p}_0 \). Hence, the computation of the boundary conjunct is numerically cheap.
2.3 A first-order formula for shape sensitivity

Taking the inner product of (9) with the adjoint solution, we get

\[ \langle \hat{p}_0^1 | \mathcal{L}_{\Omega 0,0} \hat{p}_1 \rangle + \langle \hat{p}_0^1 | \mathcal{L}_{\Omega 1,0} \hat{p}_0 \rangle \omega_1 + \langle \hat{p}_0^1 | \mathcal{L}_{\Omega 0,1} \hat{p}_0 \rangle = 0 \] (20a)

\[ \Rightarrow \left[ \mathcal{L}_{\Omega 0,0} \hat{p}_0^1 \right]_{|\theta_1} + [\hat{p}_0^1 | \mathcal{L}_\Omega] + \langle \hat{p}_0^1 | \mathcal{L}_{\Omega 1,0} \hat{p}_0 \rangle \omega_1 + \langle \hat{p}_0^1 | \mathcal{L}_{\Omega 0,1} \hat{p}_0 \rangle = 0 \] (20b)

\[- \langle \hat{p}_0^1 | r_1 \rangle_{\theta} - \omega_1 \langle \hat{p}_0^1 | D'(\omega_0) \hat{p}_0 \rangle_{\theta} + \langle \hat{p}_0^1 | \mathcal{L}_{\Omega 1,0} \hat{p}_0 \rangle \omega_1 + \langle \hat{p}_0^1 | \mathcal{L}_{\Omega 0,1} \hat{p}_0 \rangle = 0 \] (20c)

Unlike for perturbations of equation parameters, \( \hat{p}_1 \) is not completely vanishing from the equation as it still appears in the boundary conjunct \( [\hat{p}_0^1 | \mathcal{L}_\Omega] \). However, on the boundary \( \partial \Omega(0) \), the first-order drift \( \hat{p}_1 \) can be inferred from the boundary displacement – see eq. (19) – and is thus known. Hence, the above equation allows determining the eigenvalue drift at first order:

\[ \omega_1 = - \frac{\langle \hat{p}_0^1 | \mathcal{L}_{\Omega 0,1} \hat{p}_0 \rangle - \langle \hat{p}_0^1 | r_1 \rangle_{\theta}}{\langle \hat{p}_0^1 | \mathcal{L}_{\Omega 1,0} \hat{p}_0 \rangle - \langle \hat{p}_0^1 | D'(\omega_0) \hat{p}_0 \rangle_{\theta}}. \] (21)

2.4 Implementational questions

The term \( r_1 \) contains first and second order derivatives, which are to be integrated over the unperturbed domain surface. In this study, we use second-order tetrahedral finite elements, which allow computing these quantities on the triangles forming the discretized domain surface. For the Bubnov-Galerkin finite elements utilized, the discrete and the continuous adjoint formulation, i.e. first deriving the discretization and then adjoining the operators or vice versa, amount to the same discretization matrices for the adjoint equations. However, the utilized computational scheme applied to obtain the examples below is a continuous adjoint approach, because the sensitivity is computed from the continuous formulation (21). Indeed, any approach utilizing the boundary conjunct must be classified as continuous adjoint, as after the discretization the differential operators become matrices and there would not be any concept of a boundary conjunct. This is opposed to the discrete adjoint approaches using parameterized meshes to enable shape sensitivity analysis. An obvious major benefit from the continuous strategy is that there is no need to incorporate the deformations on the mesh-level, which would require specialized software, and restrict the variety of applicable mesh types. We emphasize that the Rijke tube example below uses an unstructured tetrahedral mesh, which has not been processed any further to conduct the analysis, see Fig. 2.

3 Example

As an example, we consider a tubular duct that features a zone of heat release in its interior. Such Rijke-tube configurations are commonly used as minimum working examples in the thermoacoustics community, e.g. in [1]. The Rijke tube studied here features a length of 0.5 m and a radius of \( R = 0.025 \) m. All boundaries except the outlet are considered closed (i.e. soundhard)\(^3\) the outlet is open (i.e. sound soft). The speed of sound in the first half of the tube is set to 347 m/s; 695 m/s are specified in the second half of the tube. An \( n-\tau \)-model describes the flame dynamics

\[ F(\omega) = n \exp(-i\omega \tau) \] (22)

where \( n \) is chosen to model the temperature jump for a medium with the same properties of air and \( \tau \) is 1 ms. The zone of heat release is set in the middle of the tube with the reference point being at the center axis 0.005 m upstream of the middle of the tube.

\(^3\)A sound-hard boundary condition is a simple model for sonic flow at the inlet.
Two eigenmodes, #M1 and #M2, are considered in this article; they are depicted in Fig. 3. As examples, distortions of the inlet and the outlet surface are studied. Two different distortion mappings are considered. The first one simply stretches the tube in the axial direction while the second one transforms the flat surface into a paraboloid. The two mappings are mathematically expressed as

\[ \vec{x} \mapsto \vec{x} + \epsilon \vec{R} \vec{e}_z \]  

for the shift and

\[ \vec{x} \mapsto \vec{x} + \epsilon (R - (x^2 + y^2)) \vec{e}_z \]  

for the paraboloid. The two mappings are of comparable order, as the paraboloid is designed such that a point on the center axis of the tube is displaced exactly in the same manner as when applying the planar shift.

Table 1 lists the shape sensitivities obtained from the adjoint method for various combinations of mode, distortion and considered surface. For comparison, shape sensitivities computed using a first-order forward finite difference scheme are also listed in the table. To obtain these results, new meshes were generated from the old ones by slightly displacing the grid points on the considered surfaces according to the distortion mappings (23) and (24). The finite difference step size was \( \Delta \epsilon = 1E - 9 \). Because the displacements are defined to shift the considered surfaces in the direction of the \( z \)-axis, application to the inlet and the outlet will shorten and lengthen the tube, respectively. This is in accordance with the computed sensitivities in the frequencies (real(\( \omega \))) and explains the opposite trends for the sensitivities obtained at the inlet and the outlet. Because the two displacement types are similar at first order, their sensitivities are also very similar.

However, for all reported cases the finite difference and the adjoint method show excellent agreement. The minor difference in the two methods can be attributed to a discretization error as the adjoint method discussed here follows a continuous adjoint approach. Yet, the results validate the approach.

**CONCLUSIONS**

Adjoint-based computation of the shape gradient in a Rijke tube has been presented. Comparison of the results with finite-difference-approximations show good agreement. Compared to such traditional methods the adjoint-
Figure 3. Unperturbed modes considered in the example.

Table 1. Comparison of adjoint-based results and finite difference estimates of the shape sensitivity in 1/s

<table>
<thead>
<tr>
<th>case</th>
<th>domain</th>
<th>distort</th>
<th>mode</th>
<th>adjoint</th>
<th>finite difference</th>
</tr>
</thead>
<tbody>
<tr>
<td>#T1</td>
<td>outlet</td>
<td>paraboloid</td>
<td>#M1</td>
<td>$-6.1556 - 4.054i$</td>
<td>$-6.095 - 4.060i$</td>
</tr>
<tr>
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<td>outlet</td>
<td>paraboloid</td>
<td>#M2</td>
<td>$-7.162 + 19.980i$</td>
<td>$-7.138 + 19.926i$</td>
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<td>paraboloid</td>
<td>#M1</td>
<td>$6.554 + 7.103i$</td>
<td>$6.611 + 7.199i$</td>
</tr>
<tr>
<td>#T4</td>
<td>inlet</td>
<td>paraboloid</td>
<td>#M2</td>
<td>$61.136 + 18.361i$</td>
<td>$61.904 + 18.598i$</td>
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<td>shift</td>
<td>#M1</td>
<td>$-6.236 - 4.107i$</td>
<td>$-6.195 - 4.086i$</td>
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<td>outlet</td>
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<td>$61.923 + 18.597i$</td>
<td>$62.753 + 18.888i$</td>
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</table>

based formulation requires only one solution of the direct and one solution of the adjoint problem, thereby considerably saving computational effort for the solution of nonlinear eigenvalue problems and costly remeshing.

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REFERENCES


