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Abstract

In modern science, efficient numerical treatment of high-dimensional problems becomes more and more important. A fundamental insight of the theory of information-based complexity (IBC for short) is that the computational hardness of a problem cannot be described properly only by the rate of convergence. There exist problems for which an exponential number of information operations is needed in order to reduce the initial error, although there are algorithms which provide an arbitrarily large rate of convergence. Problems that yield this exponential dependence are said to suffer from the curse of dimensionality. While analyzing numerical problems it turns out that we can often vanquish this curse by exploiting additional structural properties. The aim of this paper is to present several approaches of this type. Moreover, a detailed introduction to the field of IBC is given.

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In modern science, efficient numerical treatment of high-dimensional problems becomes more and more important. A fundamental insight of the theory of *information-based complexity* (*IBC* for short) is that the computational hardness of a problem cannot be described properly only by the rate of convergence. An impressive example that illustrates this fact was given recently by Novak and Woźniakowski [NW09]. They studied a problem for which an exponential number of information operations is needed in order to reduce the initial error, although there exist algorithms which provide an arbitrarily large rate of convergence. Problems that yield this exponential dependence are said to suffer from the *curse of dimensionality*. While analyzing numerical problems it turns out that we can often vanquish this curse by exploiting additional structural properties. The aim of this paper is to present several approaches of this type.

A numerical problem S is given by a sequence of compact linear operators S_d acting between normed spaces \mathcal{F}_d and \mathcal{G}_d , where $d \in \mathbb{N}$. In general we seek for algorithms $A_{n,d}$ that approximate S_d while using at most $n \in \mathbb{N}_0$ pieces of information on the input elements $f \in \mathcal{F}_d$. The quality of this approximation is measured by the so-called worst case error

$$\Delta^{\text{wor}}(A_{n,d}; S_d) = \sup_{\|f|\mathcal{F}_d\| \le 1} \|S_d(f) - A_{n,d}(f) \mid \mathcal{G}_d\|,$$

which we try to minimize. Problems based on tensor product structures, as well as linear algorithms that are easy to implement, are of particular interest. The minimal number of information operations needed to solve the given problem S to within a threshold $\varepsilon > 0$ is called *information complexity*:

$$n(\varepsilon,d;S_d) = \min\left\{n \in \mathbb{N}_0 \mid \exists A_{n,d} \colon \Delta^{\mathrm{wor}}(A_{n,d};S_d) \leq \varepsilon\right\}, \quad \varepsilon > 0, \, d \in \mathbb{N}.$$

If this quantity grows exponentially fast with the dimension d then S suffers from the curse of dimensionality. In the case where $n(\varepsilon, d; S_d)$ is exponential neither in d nor in ε^{-1} , the problem S is said to be weakly tractable. A special case is described by the notion of polynomial tractability, for which the information complexity needs to be bounded from above by a polynomial in d and ε^{-1} , i.e.

$$n(\varepsilon, d; S_d) \leq C \, \varepsilon^{-p} \, d^q$$
 for some $C, p > 0, \, q \geq 0$ and all $\varepsilon \in (0, 1], \, d \in \mathbb{N}$.

If the latter inequality is valid even for q=0 then S is called strongly polynomially tractable.

Next we present the three approaches to exploit structural properties we study in this paper and we briefly summarize our main complexity results.

A rather simple class of problems S is given by the set of all compact linear operators between tensor products of Hilbert spaces. Especially the complexity of tensor product problems $S_d = \bigotimes_{k=1}^d S_1 \colon H_d \to \mathcal{G}_d$, induced by some operator $S_1 \colon H_1 \to \mathcal{G}_1$, is wellunderstood. It depends on the non-increasingly ordered sequence $\lambda = (\lambda_m)_{m \in \mathbb{N}}$ of the squares of the singular values of the underlying operator S_1 . In particular, it is well-known that $S = (S_d)_{d \in \mathbb{N}}$ is not polynomially tractable if we have $\lambda_1 \geq 1$ and $\lambda_2 > 0$. Actually, we are faced with the curse of dimensionality if λ_1 is strictly larger than 1 and $\lambda_2 > 0$, or if $\lambda_1 \geq \lambda_2 = 1$ (cf. Theorem 2.11).

A first approach to modify such a problem is to scale the inner products of the source spaces H_d , $d \in \mathbb{N}$. We set

$$\langle \cdot, \cdot \rangle_{\mathcal{F}_d} = \frac{1}{s_d} \langle \cdot, \cdot \rangle_{H_d}$$
 for some $s_d > 0$ and all $d \in \mathbb{N}$

and investigate the complexity of the problem operators S_d interpreted as mappings between the Hilbert spaces \mathcal{F}_d and \mathcal{G}_d , $d \in \mathbb{N}$. The resulting problem, scaled by factors from the sequence $s=(s_d)_{d\in\mathbb{N}}$, is then denoted by $S_{(s)}=(S_{d,s_d}\colon\mathcal{F}_d\to\mathcal{G}_d)_{d\in\mathbb{N}}$. We study the worst case setting with respect to the absolute error criterion and prove

Theorem 1. With the above notation and assuming that $\lambda_2 > 0$, the following assertions are equivalent:

- (I) $S_{(s)}$ is strongly polynomially tractable.
- (II) $S_{(s)}$ is polynomially tractable.
- (III) There exists $\tau \in (0, \infty)$ such that $\lambda \in \ell_{\tau}$ and $\sup_{d \in \mathbb{N}} s_d \|\lambda\| \ell_{\tau}\|^d < \infty$.
- (IV) There exists $\varrho \in (0, \infty)$ such that $\lambda \in \ell_{\varrho}$ and $\limsup_{d \to \infty} s_d^{1/d} < 1/\lambda_1$.

If one (and hence all) of these conditions applies then the exponent of strong polynomial tractability is given by $p^* = \inf\{2\tau \mid \tau \text{ fulfills condition (III)}\}.$

We refer to Theorem 3.2 in Section 3.2.1. It is remarkable that similar to unscaled problems, polynomial tractability of the problem $S_{(s)}$ already implies strong polynomial tractability, despite the fact that we can choose the sequence of scaling factors $(s_d)_{d\in\mathbb{N}}$ completely arbitrary.

The feature that $S_{(s)}$ is weakly tractable (which is less restrictive than polynomial tractability) and the curse of dimensionality can be characterized, provided that we additionally assume a certain asymptotic behavior of the initial error $\varepsilon_d^{\rm init} = \sqrt{s_d \lambda_1^d}$ (see Theorem 3.7 in Section 3.2.2).

THEOREM 2. We study the scaled tensor product problem $S_{(s)} = (S_{d,s_d})_{d \in \mathbb{N}}$ in the worst case setting with respect to the absolute error criterion and assume $\lambda_2 > 0$.

- Let $\ln \varepsilon_d^{\mathrm{init}} \notin o(d)$ as $d \to \infty$. Then we have the curse of dimensionality. Let $\varepsilon_d^{\mathrm{init}} \in \Theta(d^{\alpha})$ as $d \to \infty$, for some $\alpha \ge 0$.
- - \circ If $\lambda_1 = \lambda_2$ then $S_{(s)}$ suffers from the curse of dimensionality.
 - \circ In the case $\lambda_1 > \lambda_2$ the problem $S_{(s)}$ is weakly tractable if and only if λ_n is $o(\ln^{-2(1+\alpha)} n)$ as $n \to \infty$.

• Let $\varepsilon_d^{\text{init}} \to 0$ as $d \to \infty$. Then we are never faced with the curse of dimensionality. Furthermore, $S_{(s)}$ is weakly tractable if and only if

(i)
$$\lambda_1 = \lambda_2$$
 and $\lambda_n \in o(\ln^{-2} n)$ as $n \to \infty$, and $\varepsilon_d^{\rm init} \in o(1/d)$ as $d \to \infty$, or (ii) $\lambda_1 > \lambda_2$ and $\lambda_n \in o(\ln^{-2} n)$ as $n \to \infty$.

(ii)
$$\lambda_1 > \lambda_2$$
 and $\lambda_n \in o(\ln^{-2} n)$ as $n \to \infty$.

Here the parameter α that controls the polynomial growth of the initial error is of particular interest. In the case where $\lambda_1 > \lambda_2$ it directly enters the characterization of weak tractability. Moreover, condition $\varepsilon_d^{\text{init}} \in o(1/d)$ as $d \to \infty$ in the third part of the theorem is quite surprising. Since for unscaled problems the initial error can only grow or decline exponentially, or it equals one in any dimension, these phenomena cannot occur in the classical theory, i.e. when $s_d = 1$ for all $d \in \mathbb{N}$.

Another approach to overcome the curse of dimensionality is related to problems defined between function spaces. Here we can make use of some a priori knowledge about the influence of certain (groups of) variables on the functions in the source space, in order to approximate them efficiently. To this end, we endow these spaces with weighted norms. During the last years especially problems on function spaces that yield a Hilbert space structure, equipped with so-called product weights, attracted a lot of attention. Problems where the source and/or target spaces are allowed to be more general Banach spaces were studied less frequently within the IBC community.

Among other things, in this paper we consider the uniform approximation problem

$$\mathrm{App} = (\mathrm{App}_d \colon F_d^{\gamma} \to \mathrm{L}_{\infty}([0,1]^d))_{d \in \mathbb{N}} \quad \text{with} \quad \mathrm{App}_d(f) = f \quad \text{for } d \in \mathbb{N}$$

defined on certain classes of smooth functions

$$F_d^{\gamma} = \{ f \colon [0,1]^d \to \mathbb{R} \mid f \in C^{\infty}([0,1]^d) \text{ with } \|f \mid F_d^{\gamma}\| < \infty \}$$

which are endowed with the weighted norms

$$||f| F_d^{\gamma}|| = \sup_{\alpha \in \mathbb{N}_a^d} \frac{1}{\gamma_{\alpha}} ||D^{\alpha} f|| L_{\infty}([0,1]^d)||.$$

Here for every $\alpha \in \mathbb{N}_0^d$, $d \in \mathbb{N}$, the product weights $\gamma_{\alpha} = \prod_{j=1}^d (\gamma_{d,j})^{\alpha_j}$ are constructed from a uniformly bounded sequence $C_{\gamma} \geq \gamma_{d,1} \geq \cdots \geq \gamma_{d,d} > 0$ of so-called generator weights. It turns out that the complexity of the approximation problem depends on certain summability properties of these generators which also play an important role when dealing with problems on product-weighted Hilbert spaces. We define the quantities

$$p(\gamma) = \inf \Big\{ \kappa > 0 \; \Big| \; \limsup_{d \to \infty} \sum_{j=1}^{d} (\gamma_{d,j})^{\kappa} < \infty \Big\},$$

$$q(\gamma) = \inf \left\{ \kappa > 0 \mid \limsup_{d \to \infty} \sum_{i=1}^{d} \frac{(\gamma_{d,j})^{\kappa}}{\ln(d+1)} < \infty \right\},$$

and prove the following

Theorem 3. For the worst case setting with respect to the absolute error criterion we have:

- If the problem App is polynomially tractable then $q(\gamma) \leq 1$. Moreover, strong polynomial tractability implies $p(\gamma) \leq 1$.
- If $q(\gamma) < 1$ or even $p(\gamma) < 1$ then App is polynomially tractable or even strongly polynomially tractable, respectively.

In fact, we prove these necessary and sufficient criteria for a whole scale of weighted Banach spaces that fulfill certain embedding conditions (see Propositions 4.6 and 4.7 for details). The source space F_d^{γ} as defined above appears as a special case within this scale. On the other hand, it generalizes a space considered by Novak and Woźniakowski [NW09]. In addition, we prove that the sufficient conditions $q(\gamma) < 1$ and $p(\gamma) < 1$ are also necessary for (strong) polynomial tractability of the L_{∞} -approximation problem defined on a certain unanchored Sobolev space \mathcal{H}_d^{γ} (cf. Theorem 4.18).

Weak tractability and the curse of dimensionality can be characterized as follows.

THEOREM 4. For App = $(App_d)_{d\in\mathbb{N}}$ the following assertions are equivalent:

- (i) The problem is weakly tractable.
- (ii) The curse of dimensionality is not present.
- (iii) For all $\kappa > 0$ we have $\lim_{d \to \infty} \frac{1}{d} \sum_{j=1}^{d} (\gamma_{d,j})^{\kappa} = 0$.
- (iv) There exists $\kappa \in (0,1)$ such that $\lim_{d\to\infty} \frac{1}{d} \sum_{j=1}^d (\gamma_{d,j})^{\kappa} = 0$.

This immediately follows from our Theorem 4.9, in which we discuss a more general situation. Note that the implication (ii) \Rightarrow (i) is not trivial. Moreover, condition (iv) is typical for problems defined on Hilbert spaces equipped with product weights.

Finally, our third approach to vanquish the curse is based on exploiting certain symmetry properties of the elements in the source space. For this purpose we again consider tensor product problems $S = (S_d \colon H_d \to \mathcal{G}_d)_{d \in \mathbb{N}}$ between Hilbert spaces. But now we restrict them to suitable subspaces which solely consist of (anti)symmetric elements. We illustrate this concept by considering the special case of problems defined between function spaces.

For $d \in \mathbb{N}$ and $I \subseteq \{1, \ldots, d\}$ let \mathcal{S}_I denote the collection of all permutations π of the coordinate set $\{1, \ldots, d\}$ that leave the complement $I^c = \{1, \ldots, d\} \setminus I$ of I fixed. Then a real-valued function $f \in H_d = H_1 \otimes \cdots \otimes H_1$ on $[0, 1]^d$ is called I-symmetric if

$$f(\boldsymbol{x}) = f(\pi(\boldsymbol{x}))$$
 for every $\boldsymbol{x} \in [0,1]^d$ and all $\pi \in \mathcal{S}_I$.

In contrast, f is called I-antisymmetric if $f(\boldsymbol{x}) = (-1)^{|\pi|} f(\pi(\boldsymbol{x}))$ for every \boldsymbol{x} and π . In what follows we denote the corresponding linear subspaces of H_d that exclusively contain symmetric or antisymmetric functions by $\mathfrak{S}_I(H_d)$ and $\mathfrak{A}_I(H_d)$, respectively. In particular, antisymmetric functions, i.e. functions that change their sign when we exchange the variables x_i and x_j , $i,j \in I$, turned out to be of some practical interest; see, e.g., Section 5.4.2. For the restriction of a given tensor product problem $S = (S_d \colon H_d \to \mathcal{G}_d)_{d \in \mathbb{N}}$ to the subspaces $P_{I_d}(H_d)$, $d \in \mathbb{N}$, we write $S_I = (S_{d,I_d})_{d \in \mathbb{N}}$. Here the kind of symmetry $P \in \{\mathfrak{S}, \mathfrak{A}\}$, as well as a sequence $(I_d)_{d \in \mathbb{N}}$ of subsets of the coordinates, are assumed to be fixed.

Since for $d \in \mathbb{N}$ the operators S_{d,I_d} can be interpreted as a composition of S_d with suitable orthogonal projections, there exists a close relation of the singular values of S_d to the corresponding singular values of the restricted operators S_{d,I_d} . These numbers

essentially determine the minimal worst case error of the problem S_I . This knowledge furthermore allows the construction of an optimal (linear) algorithm that realizes this error (cf. Theorem 5.4).

Consequently, we can conclude assertions that relate the information complexity of S_I to the squares of the singular values of S_1 and to the number of (anti)symmetry conditions we impose. For the sake of simplicity we restrict ourselves again to the absolute error criterion and start by discussing the case of symmetric problems (see Theorem 5.10).

THEOREM 5 (Polynomial tractability, $P = \mathfrak{S}$). Let $S_1 : H_1 \to \mathcal{G}_1$ denote a compact linear operator between Hilbert spaces and let $\lambda = (\lambda_m)_{m \in \mathbb{N}}$ be the sequence of eigenvalues of $W_1 = S_1^{\dagger} S_1$ with non-increasing ordering. Assume $\lambda_2 > 0$ and for d > 1 let $\emptyset \neq I_d \subseteq \{1, \ldots, d\}$ be fixed. We consider the restriction $S_I = (S_{d,I_d})_{d \in \mathbb{N}}$ of the tensor product problem $S = (S_d : H_d \to \mathcal{G}_d)_{d \in \mathbb{N}}$ to the I_d -symmetric subspaces $\mathfrak{S}_{I_d}(H_d) \subset H_d$, $d \in \mathbb{N}$. Then S_I is strongly polynomially tractable if and only if $\lambda \in \ell_\tau$ for some $\tau \in (0, \infty)$ and

- $\lambda_1 < 1$, or
- $1 = \lambda_1 > \lambda_2$ and $d \#I_d \in \mathcal{O}(1)$ as $d \to \infty$.

Moreover, provided that $\lambda_1 \leq 1$, the problem is polynomially tractable if and only if $\lambda \in \ell_\tau$ for some $\tau \in (0, \infty)$ and

- $\lambda_1 < 1$, or
- $\lambda_1 = 1$ and $d \#I_d \in \mathcal{O}(\ln d)$ as $d \to \infty$.

The problem of finding conditions which are sufficient for polynomial tractability in the case $\lambda_1 > 1$ remains open. However, our results show that the conditions $\lambda \in \ell_{\tau}$ and $d - \#I_d \in \mathcal{O}(\ln d)$ are necessary in this situation, too. In conclusion, we see that imposing sufficiently many additional symmetry assumptions, we can avoid the curse of dimensionality which we are faced with e.g. in the case $\lambda_1 = \lambda_2 = 1$; see also Theorem 2.11.

The complexity analysis of antisymmetric problems is more demanding. On the other hand, it turns out that here even weaker conditions are sufficient to conclude polynomial tractability and thus to vanquish the curse. One of the reasons is the structure of the initial error which is more complicated in this case. Similar to Theorem 5.16 in Section 5.3 we can summarize the main results on the complexity as follows:

THEOREM 6 (Polynomial tractability, $P = \mathfrak{A}$). Let $S_1 \colon H_1 \to \mathcal{G}_1$ denote a compact linear operator between Hilbert spaces and let $\lambda = (\lambda_m)_{m \in \mathbb{N}}$ be the sequence of eigenvalues of $W_1 = S_1^{\dagger} S_1$ with non-increasing ordering. Assume $\lambda_2 > 0$ and for d > 1 let $\emptyset \neq I_d \subseteq \{1, \ldots, d\}$ be fixed. We consider the restriction $S_I = (S_{d,I_d})_{d \in \mathbb{N}}$ of the tensor product problem $S = (S_d \colon H_d \to \mathcal{G}_d)_{d \in \mathbb{N}}$ to the I_d -antisymmetric subspaces $\mathfrak{A}_{I_d}(H_d) \subset H_d$, $d \in \mathbb{N}$. Then for the case $\lambda_1 < 1$ the following statements are equivalent:

- S_I is strongly polynomially tractable.
- S_I is polynomially tractable.
- There exists a constant $\tau \in (0, \infty)$ such that $\lambda \in \ell_{\tau}$.

Moreover, the same equivalences hold true if $\lambda_1 \geq 1$ and the number of antisymmetric coordinates $\#I_d$ grows linearly with the dimension d.

Clearly, these assertions show that antisymmetric tensor product problems are significantly easier than their symmetric counterparts, which on their part possess a lower information complexity than entire tensor product problems, as long as we impose enough (anti)symmetry conditions. On the other hand, there exist quite natural examples which show that even fully antisymmetric problems are not necessarily trivial or polynomially tractable in general. For details we refer the reader to Section 5.4.1.

Let us briefly explain the structure of the present paper. In the first chapter we settle some notational conventions and define the abstract problem we are faced with in IBC. Furthermore, the cost model we are going to use is introduced and we recall the formal definitions of several complexity categories.

In Chapter 2 we discuss special classes of numerical problems, as well as elementary tools that we need to handle them. In particular, we give a detailed introduction to the singular value decomposition (SVD) of compact operators between Hilbert spaces. In many cases it forms the basis for the construction of optimal algorithms. Hence it is of fundamental importance for the rest of our work. In addition, we discuss tensor product structures in Hilbert spaces and recall some well-known complexity assertions for problems related to this concept. Finally, we briefly introduce so-called reproducing kernel Hilbert spaces (RKHSs) and collect some of their properties.

In the first two sections of the third chapter we derive characterizations of the different types of tractability of scaled tensor product problems between Hilbert spaces we presented in Theorems 1 and 2 above. Moreover, from them we derive a complete characterization for the normalized error criterion in Section 3.2.3. It turns out that here the scaling factors become irrelevant. Apart from formulas of the optimal algorithm and its worst case error, we additionally show that these new assertions generalize the known theory in a quite natural way. We conclude this chapter by the application of the results obtained to two simple examples.

Chapter 4 deals with problems on function spaces endowed with weighted norms. Here we explain the concept of weighted spaces in full detail and illustrate it using the example of an unanchored Sobolev \mathcal{H}_d^{γ} space equipped with product weights. For the uniform approximation problem on this space we present an algorithm $A_{n,d}^*$ that satisfies suitable upper error bounds. Together with corresponding lower bounds, which we prove for spaces of low-degree polynomials, the application of simple embedding arguments then leads us to complexity assertions for a whole scale of product-weighted Banach spaces. In particular, these assertions cover the results for the space F_d^{γ} stated in Theorems 3 and 4. Finally, the last section of the fourth chapter, Section 4.4, presents some generalizations of the techniques developed before. Among other things, we show how to handle L_p -approximation problems, where $1 \le p < \infty$, defined on suitable spaces. Moreover, we show that the algorithm $A_{n,d}^*$ is essentially optimal for L_{∞} -approximation on \mathcal{H}_d^{γ} . For the proof we make use of arguments due to Kuo, Wasilkowski and Woźniakowski [KWW08] that relate the uniform approximation problem in the worst case setting, defined on quite general reproducing kernel Hilbert spaces, to a certain average case L₂-approximation problem.

Some of the results presented in Chapter 4 were already published in [Wei12b]. However, we were able to partially improve them. We explicitly emphasize generalizations and new results at appropriate points.

Finally, Chapter 5 is devoted to problems with (anti)symmetry conditions. We start with the definition of (anti)symmetry in Hilbert function spaces. In particular, we focus on tensor product structures and deduce fundamental properties of the respective projections and subspaces. At the end of Section 5.1 we use these properties to generalize the notion of (anti)symmetry to tensor products of abstract Hilbert spaces. Afterwards we define (anti)symmetric numerical problems $S_I = (S_{d,I_d})_{d \in \mathbb{N}}$ by the restriction of a given tensor product problem $S = (S_d)_{d \in \mathbb{N}}$ to the subspaces of (anti)symmetric elements in the source spaces. We prove the commutativity of the operators S_d with certain projections and deduce formulas for optimal algorithms and their worst case errors. In Section 5.3 we then discuss the complexity of (anti)symmetric numerical problems. We distinguish between symmetric and antisymmetric problems, as well as between the absolute and the normalized error criterions. Here we in particular derive the proofs of Theorems 5 and 6. The chapter is concluded by a section devoted to several applications. On the one hand, we use simple examples to show that the additional knowledge about (anti)symmetry conditions can dramatically reduce information complexity. On the other hand, we also discuss more advanced problems that play a role in computational practice. To this end, we illustrate the application of this new theory to the approximation problem of socalled wavefunctions that arise in certain models of quantum mechanics and theoretical chemistry.

A major part of the results proven in Chapter 5 was published in [Wei12a]. However, at some points we use different proof techniques that allow slight generalizations.

The symbols \square and \blacksquare are used to indicate the end of remarks and examples, and of proofs, respectively.

1. Preliminaries

Apart from introducing some notational conventions, the aim of this chapter is to define the general objects of interest in *information-based complexity*. We give an abstract formulation of the general problem in Section 1.2. Afterwards, in Section 1.3, we introduce some classes of algorithms and discuss the cost model we are going to use in this paper. Finally, in Section 1.4, we recall the notions of tractability, as well as the definition of the curse of dimensionality.

1.1. Basic notation. As usual, we denote by \mathbb{N} the natural numbers, and $\mathbb{N}_0 = \mathbb{N} \cup \{0\}$ are all non-negative integers. Moreover, \mathbb{R} denotes the real line and \mathbb{R}^d $(d \in \mathbb{N})$ is the collection of all points $\boldsymbol{x} = (x_1, \dots, x_d)$ in the d-dimensional Euclidean space. Given a real number y > 0, the symbol $\lfloor y \rfloor$ means the largest $n \in \mathbb{N}_0$ such that $n \leq y$, and we define $\lceil y \rceil$ to be the smallest number $m \in \mathbb{N}$ with $y \leq m$. The value of the Riemann zeta function at some z > 1 is denoted by $\zeta(z) = \sum_{n=1}^{\infty} n^{-z}$.

If $\mathbf{k} = (k_1, \dots, k_d) \in \mathbb{N}_0^d$ is a multi-index then $|\mathbf{k}| = \sum_{i=1}^d k_i$ stands for its length. Furthermore, we use the common notation $\mathbf{x}^{\mathbf{k}} = x_1^{k_1} \cdot \dots \cdot x_d^{k_d}$. For $\alpha \in \mathbb{N}_0^d$ partial derivatives of d-variate functions are denoted by D^{α} , i.e.

$$D^{\alpha} f = \frac{\partial^{|\alpha|} f}{\partial x_1^{\alpha_1} \cdots \partial x_d^{\alpha_d}}.$$

Derivatives of univariate functions g are indicated by $g', g'', \ldots, g^{(n)}$, respectively. For real numbers a < b half-open intervals are denoted by [a, b), and $[a, b]^d$ stands for the Cartesian product $\times_{i=1}^d [a, b] = [a, b] \times \ldots \times [a, b]$. If $\boldsymbol{x} \in \mathbb{R}^d$ belongs to $[a, b]^d$ then the value of the characteristic (or indicator) function $\chi_{[a,b]^d}(\boldsymbol{x})$ of this set equals 1. Otherwise we define $\chi_{[a,b]^d}(\boldsymbol{x}) = 0$. Similarly the Kronecker delta function $\delta_{i,j}$ is one if i and j coincide and $\delta_{i,j} = 0$ otherwise.

We assume that the reader has a fundamental knowledge in measure theory and probability theory as can be found, e.g., in the textbooks of Bauer [Bau01, Bau96]. We write λ^d for the Lebesgue measure in \mathbb{R}^d and use the symbols \mathbb{P} and \mathbb{E} for probabilities and expectations, respectively. We use #I to denote the cardinality of a finite set I. As usual, the sum over an empty index set I is to be interpreted as zero, whereas empty products equal 1 by definition.

We also assume that the reader is familiar with the basic concepts in functional analysis such as, e.g., complete normed spaces (Banach spaces), weak derivatives or tensor products. For a comprehensive introduction we refer to the textbooks of Triebel [Tri92] and Yosida [Yos80]. The norm in some space F is denoted by $\|\cdot\|F\|$. We write $B_r(F) = 0$

 $\{f \in F \mid \|f \mid F\| \leq r\}$ for closed balls of radius $r \geq 0$ centered at 0 in normed spaces F. Moreover, we use ∂M for the boundary and $\operatorname{int}(M)$ for the interior of a set M. Furthermore, $\mathcal{B}(F) = B_1(F) = \operatorname{int}(B_1(F)) \cup \partial B_1(F)$ denotes the unit ball in F. For the class of all bounded linear operators between normed spaces F and G we write $\mathcal{L}(F,G)$. The subset of all compact operators is denoted by $\mathcal{K}(F,G)$. We say a space F is (continuously) embedded into another space G with norm G if the operator norm of id: $F \to G$, $f \mapsto \operatorname{id}(f) = f$, equals $G \in [0, \infty)$. In this case we write $F \hookrightarrow G$ and $\|\operatorname{id} \mid \mathcal{L}(F,G)\| = C$. We use $\langle \cdot, \cdot \rangle_H$ for the inner product in a Hilbert space H. Moreover, we write M^{\perp} for the orthogonal complement of a linear subspace $M \subset H$ and we use \oplus to denote the orthogonal sum with respect to $\langle \cdot, \cdot \rangle_H$.

If $(\mathcal{X}, \mathfrak{a}, \mu)$ is an arbitrary measure space and $0 then we use <math>L_p(\mathcal{X}, \mathfrak{a}, \mu)$ to denote the classical Lebesgue spaces. Hence, if $p < \infty$ then we deal with the set of (equivalence classes of) μ -measurable functions $f \colon \mathcal{X} \to \mathbb{R}$ for which the norm (1)

$$||f| L_p(\mathcal{X}, \mathfrak{a}, \mu)|| = \left(\int_{\mathcal{X}} |f(x)|^p d\mu(x)\right)^{1/p}$$

is finite. Moreover, $L_{\infty}(\mathcal{X}, \mathfrak{a}, \mu)$ is the space (of classes) of μ -essentially bounded functions on \mathcal{X} , furnished with the norm

$$||f| L_{\infty}(\mathcal{X}, \mathfrak{a}, \mu)|| = \operatorname{ess\,sup}_{x \in \mathcal{X}} |f(x)|.$$

As usual, two functions are identified if they coincide μ -almost everywhere on \mathcal{X} , and we do not distinguish between functions and their equivalence classes. The following special cases are of particular interest for us.

For a Borel measurable subset $\mathcal{X} = \Omega \subset \mathbb{R}^d$, the Borel sigma-algebra $\mathfrak{a} = \Sigma = \Sigma(\Omega)$, and $\mu = \lambda^d$ we use the shorthand $L_p(\Omega) = L_p(\Omega, \Sigma, \lambda^d)$. If μ is not the Lebesgue measure, but is absolutely continuous with respect to λ^d , and if $\varrho = \mathrm{d}\mu/\mathrm{d}\lambda^d$ describes a probability density function that is strictly positive $(\lambda^d$ -a.e.) on Ω , then we write $L_p^\varrho(\Omega)$. On the other hand, for a discrete measure space $(\Gamma, \mathfrak{b}, \nu)$ on some set Γ with $\nu(\{i\}) = 1$ for each $i \in \Gamma$ we write $\ell_p(\Gamma) = L_p(\Gamma, \mathfrak{b}, \nu)$ and we abbreviate the notation to ℓ_p if $\Gamma = \mathbb{N}$. Keep in mind that in this case the norms simplify to

$$\|\lambda \mid \ell_p\| = \begin{cases} \left(\sum_{m=1}^{\infty} |\lambda_m|^p\right)^{1/p} & \text{if } 0$$

where $\lambda = (\lambda_m)_{m \in \mathbb{N}}$ is any real-valued sequence such that the above norm is finite.

Finally, we make use of the Bachmann–Landau notation for asymptotic growth rates. That is, for real-valued functions f and g defined on some subset of the real line we write $f(x) \in \mathcal{O}(g(x))$ as $x \to a$ if there exists a universal constant M > 0 such that

$$|f(x)| \le M \, |g(x)|$$

for all x sufficiently close to a. If g is non-zero (at least in a neighborhood of a) then this

⁽¹⁾ Actually, in the case 0 the given formula only provides a quasi-norm, i.e. then we need an additional constant <math>k > 1 for the triangle inequality. Since this does not play any role in our applications we do not emphasize this difference in what follows.

definition equivalently reads

$$\limsup_{x \to a} \left| \frac{f(x)}{g(x)} \right| < \infty.$$

If we have $f(x) \in \mathcal{O}(g(x))$ and simultaneously $g(x) \in \mathcal{O}(f(x))$ as $x \to a$, then we write $f(x) \in \Theta(g(x))$, $x \to a$. Moreover, we say that $f(x) \in o(g(x))$ as $x \to a$ if for any $\delta > 0$ there exists a neighborhood U of a such that

$$|f(x)| \le \delta |g(x)|$$

for all $x \in U$. Again for non-vanishing q this property can be reformulated as

$$\lim_{x \to a} \left| \frac{f(x)}{g(x)} \right| = 0.$$

All these three notations will be used in particular for sequences $(f_d)_{d\in\mathbb{N}}$ (interpreted as special classes of functions), where we have $a=\infty$.

1.2. General problem. In numerous applications in physics, chemistry, finance, economics, and computer science we are faced with very high-dimensional continuous problems which can almost never be solved analytically. Therefore we search for algorithms which approximate the unknown solutions numerically to within a threshold $\varepsilon > 0$.

In general, such a problem is given by a non-trivial solution operator

$$S \colon \widetilde{\mathcal{F}} \to \mathcal{G},$$
 (1.1)

mapping a problem element f from a subset $\widetilde{\mathcal{F}}$ of some normed space \mathcal{F} onto its solution S(f) in some (other) target space \mathcal{G} . Often, but not always, $\widetilde{\mathcal{F}}$ is assumed to be the unit ball $\mathcal{B}(\mathcal{F})$ in some Banach space \mathcal{F} of multivariate functions $f \colon \Omega_d \subset \mathbb{R}^d \to \mathbb{R}$. For the domain of definition Ω_d , usually the unit cube $[0,1]^d$ is taken. Since the dependence on d will play a crucial role in this paper, we concentrate on whole sequences $S = (S_d)_{d \in \mathbb{N}}$ of solution operators, where every

$$S_d : \widetilde{\mathcal{F}}_d \to \mathcal{G}_d, \quad d \in \mathbb{N},$$
 (1.2)

is of the form (1.1).

Typically, $\widetilde{\mathcal{F}}_d$ is an infinite-dimensional subset of the *source space* \mathcal{F}_d and thus we cannot input $f \in \widetilde{\mathcal{F}}_d$ directly into the computer. Instead we assume that the input for our algorithms A consists of finitely many cleverly chosen pieces of information which hopefully describe f as well as possible. In Section 1.3 we define different kinds of information operations which lead us to different classes \mathcal{A}_d of algorithms. For now assume $A_d \colon \widetilde{\mathcal{F}}_d \to \mathcal{G}_d$ is a fixed element in some class \mathcal{A}_d .

The local error $\Delta_{\text{loc}}(f; A_d, S_d)$ of a given algorithm $A_d \in \mathcal{A}_d$ applied to a problem element $f \in \widetilde{\mathcal{F}}_d$ is defined as the difference of the exact solution $S_d(f)$ and the approximate solution $A_d(f)$, measured in the norm of the target space \mathcal{G}_d , i.e.

$$\Delta_{\text{loc}}(f; A_d, S_d) = ||S_d(f) - A_d(f)||\mathcal{G}_d||.$$

Based on this definition there are several ways to quantify the quality of A_d .

In the worst case setting this is done in terms of the maximal local error of the algorithm among all possible inputs $f \in \widetilde{\mathcal{F}}_d$. Hence, we define by

$$\Delta^{\text{wor}}(A_d; S_d \colon \widetilde{\mathcal{F}}_d \to \mathcal{G}_d) = \sup_{f \in \widetilde{\mathcal{F}}_d} \Delta_{\text{loc}}(f; A_d, S_d)$$

the worst case error of the algorithm A_d for the problem $S_d \colon \widetilde{\mathcal{F}}_d \to \mathcal{G}_d$. On the other hand, sometimes it is useful to measure the average performance of a given algorithm on the input set $\widetilde{\mathcal{F}}_d$. This corresponds to the so-called average case setting. Here we need to assume in addition that $\widetilde{\mathcal{F}}_d$ is equipped with a probability measure μ_d . The term

$$\Delta^{\operatorname{avg}}(A_d; S_d \colon \widetilde{\mathcal{F}}_d \to \mathcal{G}_d) = \left(\int_{\widetilde{\mathcal{F}}_d} \Delta_{\operatorname{loc}}(f; A_d, S_d)^2 \, \mathrm{d}\mu_d(f) \right)^{1/2}$$

then denotes the average case error of A_d (2). Since the worst case setting seems to be much more important we will mainly deal with worst case errors. However, for some problems there exist close relations to the average case setting. One such example will be presented in Section 4.4.2. For the sake of completeness we stress that there exist even more settings which are subject to current research; we mention the probabilistic and the randomized setting and refer to [NW08, Section 3.2] for an extensive discussion.

In numerical analysis, a major assumption is that information is expensive. Therefore we are interested in algorithms which solve a given problem within a tolerance ε while using as few pieces of information on the inputs as possible. This property can be captured by the concept of the *nth minimal error*

$$e^{\mathrm{sett}}(n, d; S_d \colon \widetilde{\mathcal{F}}_d \to \mathcal{G}_d) = \inf_{A_{n,d} \in \mathcal{A}_d^n} \Delta^{\mathrm{sett}}(A_{n,d}; S_d \colon \widetilde{\mathcal{F}}_d \to \mathcal{G}_d)$$

for sett $\in \{\text{wor}, \text{avg}\}$, $d \in \mathbb{N}$, and $n \in \mathbb{N}_0$, where the infimum is taken over all algorithms in the class

 $\mathcal{A}_d^n = \{ A \in \mathcal{A}_d \mid A \text{ uses at most } n \text{ information operations on the input} \}.$

Consequently, the initial error

$$\varepsilon_d^{\text{init,sett}} = e^{\text{sett}}(0, d; S_d : \widetilde{\mathcal{F}}_d \to \mathcal{G}_d), \quad d \in \mathbb{N},$$

describes the smallest error we can achieve without using any information on the input in a given setting sett $\in \{\text{wor}, \text{avg}\}$. We will see in Section 2.2 that under mild assumptions this initial error can be attained by the zero algorithm, i.e. by $A_{0,d} \equiv 0 \in \mathcal{G}_d$.

If there is no danger of confusion we abbreviate the above notations and simply write $\Delta^{\text{sett}}(A_d; S_d)$ and $e^{\text{sett}}(n, d; S_d)$, where sett is an element of {wor, avg}, or even only $\Delta(A_d)$ and e(n, d), respectively. Moreover, in Chapters 4 and 5 it seems to be useful to stress the source spaces \mathcal{F}_d the problem elements come from, rather than the operator S_d . There we slightly abuse notation and write $e^{\text{wor}}(n, d; \mathcal{F}_d)$ instead of $e^{\text{wor}}(n, d; S_d : \mathcal{B}(\mathcal{F}_d) \to \mathcal{G}_d)$.

The main goal in the classical theory is to find sharp bounds on the nth minimal error in terms of the amount of information operations. In fact, there is a huge literature where

⁽²⁾ In fact, μ_d is defined on the Borel sets of $\widetilde{\mathcal{F}}_d$ and we need to claim $\Delta_{\text{loc}}(\cdot; A_d, S_d)$ to be a measurable function, but these are only formal issues. See, e.g., [NW08, p. 129] for further details.

the existence of constants c_d , $C_d > 0$ and p_d , $P_d > 0$ was proven such that estimates of the type

 $\frac{1}{c_d} n^{-p_d} \le e(n, d) \le C_d n^{-P_d} \quad \text{ for all } n \in \mathbb{N}$

hold for certain problems S in a given setting (3). Back then, the respective researchers did not pay much attention to the constants c_d and C_d involved. These numbers can be arbitrarily large and in some cases their dependence on d is completely unknown. Instead the attention was focused on the so-called rate (or order) of convergence, i.e. on proofs which yield $p_d = P_d$. Often this rate tends to zero as $d \to \infty$. Therefore such bounds are not meaningful at all for large d. Thus, usually the parameter d was assumed to be a fixed (and reasonably small) constant in this approach. Since we also want to work in huge dimensions, a more careful error analysis is needed.

1.3. Algorithms and a cost model. For fixed $n, d \in \mathbb{N}$ an algorithm $A_{n,d} \in \mathcal{A}_d^n$ is modeled as a mapping $\varphi_n \colon \mathbb{R}^n \to \mathcal{G}_d$ and a function $N_n \colon \widetilde{\mathcal{F}}_d \to \mathbb{R}^n$ such that $A_{n,d} = \varphi_n \circ N_n$. For the sake of completeness in the case n = 0 we simply assign a constant value $c \in \mathcal{G}_d$ to every element $f \in \widetilde{\mathcal{F}}$, i.e. $A_{0,d} \equiv c$, in order to model an algorithm that does not depend on the input at all. If n > 0 then the information map N_n is given by

$$N_n(f) = (L_1(f), \dots, L_n(f)), \quad f \in \widetilde{\mathcal{F}}_d, \tag{1.3}$$

where $L_j \in \Lambda$. Here we distinguish certain classes of information operations Λ . In one case we assume that we are allowed to compute arbitrary continuous linear functionals on the inputs f. Then $\Lambda = \Lambda^{\rm all}$ coincides with \mathcal{F}_d^* , the dual space of \mathcal{F}_d . If we deal with problem operators S_d defined on function spaces $\widetilde{\mathcal{F}}_d$ then often only function evaluations are permitted, i.e. $L_j(f) = f(t^{(j)})$ for a certain fixed $t^{(j)} \in \Omega_d$ in the domain of definition of f. In this case $\Lambda = \Lambda^{\rm std}$ is called standard information. If function evaluation is continuous for all $t \in \Omega_d$ we have $\Lambda^{\rm std} \subset \Lambda^{\rm all}$. In particular this is the case when dealing with problems defined on reproducing kernel Hilbert spaces (RKHS) (see Section 2.5). If L_j depends continuously on f but is not necessarily linear then the respective class is denoted by $\Lambda^{\rm cont}$. Note that in this case also N_n is continuous and we obviously have $\Lambda^{\rm all} \subset \Lambda^{\rm cont}$.

Furthermore, we distinguish between adaptive and non-adaptive algorithms. The latter case is described above in formula (1.3), where L_j does not depend on the previously computed values $L_1(f), \ldots, L_{j-1}(f)$. In contrast, we also discuss algorithms of the form $A_{n,d} = \varphi_n \circ N_n$ with

$$N_n(f) = (L_1(f), L_2(f; y_1), \dots, L_n(f; y_1, \dots, y_{n-1})), \quad f \in \widetilde{\mathcal{F}}_d,$$
 (1.4)

where $y_1 = L_1(f)$ and $y_j = L_j(f; y_1, \ldots, y_{j-1})$ for $j = 2, \ldots, n$. If N_n is adaptive we restrict ourselves to the case where L_j depends linearly on f, e.g. $L_j(\cdot; y_1, \ldots, y_{j-1})$ is in $\Lambda^{\rm all}$. Note that in any case N_n is either continuous, or constructed out of linear information operations (which may be combined adaptively). Moreover, in all cases of information maps, the mapping φ_n can be chosen arbitrarily.

 $^(^3)$ In many cases those estimates hold modulo $\log n$ to some power which usually depends linearly on d. For simplicity we omit these factors because they are not crucial for the following argument.

For upper error bounds small classes of algorithms are most important. The smallest such class is the family of linear, non-adaptive algorithms of the form

$$A_{n,d}(f) = \sum_{j=1}^{n} L_j(f)g_j$$
 (1.5)

with some $g_j \in \mathcal{G}_d$ and $L_j \in \Lambda^{\operatorname{all}}$ or even $L_j \in \Lambda^{\operatorname{std}}$. We denote this set of algorithms by $\mathcal{A}_d^{n,\operatorname{lin}}(\Lambda)$, where $\Lambda = \Lambda^{\operatorname{all}}$ or $\Lambda = \Lambda^{\operatorname{std}}$, respectively. On the other hand, it is reasonable to prove lower error bounds for preferably large classes of algorithms. The most general families consist of algorithms $A_{n,d} = \varphi_n \circ N_n$, where φ_n is completely arbitrary and N_n uses either non-adaptive continuous or adaptive linear information. We denote the respective classes by $\mathcal{A}_d^{n,\operatorname{cont}}$ and $\mathcal{A}_d^{n,\operatorname{adapt}}$.

One of the most fundamental assumptions in IBC is that we can perform (exact) basic arithmetic operations on elements of the target space \mathcal{G}_d , as well as on real numbers, with unit cost. Formally this means that we work with the real number model, in contrast to the bit number model, which is used in some other fields of computational science (see, e.g., [NW08, Section 4.1.3]). Moreover, we assume that information operations on the input are given by certain black box computations which are sometimes called oracle calls. Typically the computational costs for information operations are much higher than for simple arithmetic operations since the computation of a function value or a linear functional may require billions of such operations. If we assume that every oracle call has a fixed cost $C \gg 1$ then the total cost of computing the output of an algorithm is proportional to the number of needed information operations (4). Therefore it is reasonable to study not only the nth minimal error of a given problem but also the inverse quantity which we call information complexity:

$$n_{\text{abs}}^{\text{sett}}(\varepsilon, d; S_d \colon \widetilde{\mathcal{F}}_d \to \mathcal{G}_d) = \min\{n \in \mathbb{N}_0 \mid \exists A \in \mathcal{A}_d^n \text{ such that } \Delta^{\text{sett}}(A) \le \varepsilon\}$$
$$= \min\{n \in \mathbb{N}_0 \mid e^{\text{sett}}(n, d) \le \varepsilon\},$$

where $d \in \mathbb{N}$, $\varepsilon > 0$ and sett $\in \{\text{wor, avg}\}$. That is, we look at the amount of oracle calls needed to compute an ε -approximation in dimension d. Hence, due to our assumptions this information complexity roughly equals the *total complexity* of a given problem and therefore describes its computational hardness. For a detailed discussion of algorithms and their costs, as well as the relations of information complexity and total complexity, we refer the reader to Section 4.1 in [NW08].

Finally we want to mention that the above definition addresses the absolute error criterion. In contrast we will also consider the normalized error criterion where we seek for the minimal number of information operations needed to improve the initial error by some factor $\varepsilon' > 0$. We denote the corresponding information complexity by

$$n_{\text{norm}}^{\text{sett}}(\varepsilon', d; S_d : \widetilde{\mathcal{F}}_d \to \mathcal{G}_d) = n_{\text{abs}}^{\text{sett}}(\varepsilon' \varepsilon_d^{\text{init}}, d; S_d : \widetilde{\mathcal{F}}_d \to \mathcal{G}_d), \quad \text{sett} \in \{\text{wor, avg}\}.$$

Obviously the two notions coincide if the problem under consideration is well-scaled, that is, $\varepsilon_d^{\rm init}=1$. Otherwise the problem may be significantly harder with respect to

⁽⁴⁾ There also exist approaches in which the cost of an oracle call depends on the parameters of the problem. Those attempts stress that the computational effort for function evaluations increases with the number of (active) variables. See, e.g., [KSWW10a] for details.

the normalized error criterion, e.g. if $\varepsilon_d^{\rm init}$ is exponentially small in d. Of course also the converse situation is conceivable. However, note that both the information complexities are always non-increasing in the first argument and we have

$$n_{\text{norm}}^{\text{sett}}(1, d; S_d : \widetilde{\mathcal{F}}_d \to \mathcal{G}_d) = n_{\text{abs}}^{\text{sett}}(\varepsilon_d^{\text{init}}, d; S_d : \widetilde{\mathcal{F}}_d \to \mathcal{G}_d) = 0$$
 (1.6)

for all $d \in \mathbb{N}$ due to the definition of the initial error.

Again we will use shorthands such as $n_{\text{abs}}(\varepsilon, d; S_d)$ or even $n(\varepsilon, d)$ to simplify notation.

1.4. Notions of tractability. As already indicated we strongly believe that it is not sufficient to study only the rate of convergence, i.e. the dependence of $n(\varepsilon, d)$ on ε , to properly describe the computational hardness of a given problem. We also need to incorporate the dependence on the parameter d. Keep in mind that the following definitions equally refer to both, the absolute and the normalized, error criteria. Therefore we simply write $n(\varepsilon, d)$ instead of $n_{\text{abs}}(\varepsilon, d)$ or $n_{\text{norm}}(\varepsilon, d)$ for the information complexity.

When dealing with multivariate problems we often observe the so-called curse of dimensionality, which goes back to Bellman in the late 1950s (cf. [Bel57]). Given a concrete setting, a problem is said to suffer from the curse of dimensionality if the corresponding information complexity $n(\varepsilon, d)$ increases exponentially with the dimension d. That is, for at least one $\varepsilon > 0$ there exist positive constants C and γ which are independent of the dimension such that

$$n(\varepsilon, d) \ge C(1 + \gamma)^d$$

for infinitely many $d \in \mathbb{N}$. More generally, if $n(\varepsilon, d)$ depends exponentially on d or ε^{-1} then the problem is called *intractable* (5). Otherwise we have *tractability*, which goes back to Woźniakowski in the early 1990s (see [Woź94a, Woź94b]). At that time a problem was called tractable if its complexity depends at most polynomially on ε^{-1} and d. Today this is just one case in a whole hierarchy of notions of tractability. We describe these classes starting with the weakest notion.

If a problem is not intractable then we have weak tractability, which can be equivalently expressed by

$$\lim_{\varepsilon^{-1}+d\to\infty} \frac{\ln n(\varepsilon,d)}{\varepsilon^{-1}+d} = 0$$

(see [GW08, NW08]). Here the limit is taken with respect to all two-dimensional sequences $((\varepsilon_k, d_k))_{k \in \mathbb{N}} \subset (0, 1] \times \mathbb{N}$ such that $\varepsilon_k < \varepsilon_{d_k}^{\text{init}}$ and $\varepsilon_k^{-1} + d_k \to \infty$ as $k \to \infty$. In particular, the latter restriction ensures that $n(\varepsilon, d) \geq 1$. Furthermore, we want to stress that weak tractability implies the absence of the curse of dimensionality, but in general the converse is not true. Recently a slightly stronger notion called *uniform weak tractability* has been suggested. We will not follow this line of research and refer the reader to [Sie13].

⁽⁵⁾ Formally that means that there exist universal constants $\gamma, C > 0$, as well as sequences $(\varepsilon_k)_{k \in \mathbb{N}}$ and $(d_k)_{k \in \mathbb{N}}$ with $\varepsilon_k \in (0,1]$ and $d_k \in \mathbb{N}$ for all $k \in \mathbb{N}$, such that $\varepsilon_k^{-1} + d_k \to \infty$ as $k \to \infty$, and $n(\varepsilon_k, d_k) \geq C(1+\gamma)^{\varepsilon_k^{-1} + d_k}$ for every $k \in \mathbb{N}$. Note that this definition includes the curse as a special case, where $\varepsilon_k \equiv \varepsilon_0$.

Since there are many ways to measure the lack of exponential dependence, the abstract notion of generalized (or T-) tractability was introduced (see [GW07, GW09] and [NW08, Chapter 8]). Here the essence is to describe the behavior of the information complexity in terms of a multiple of some power of a so-called tractability function T depending on ε^{-1} and d. Without going into details we mention that the following classes can be seen as special cases in this general framework.

For the sake of completeness we also introduce the notion of quasi-polynomial tractability, which was developed quite recently. A problem is called quasi-polynomially tractable if there are universal constants C, t > 0 such that

$$n(\varepsilon, d) \le C \exp(t(1 + \ln \varepsilon^{-1})(1 + \ln d))$$

for every $\varepsilon \in (0,1]$ and $d \in \mathbb{N}$. Note that for fixed ε or d this upper bound behaves polynomially in the other argument, which somehow justifies the name of this class of problems. For details see [GW11] and [NW12].

Finally, the most important and until now most studied type of tractability is called polynomial tractability. We say that a problem is polynomially tractable if there exist absolute constants C, p > 0 and $q \ge 0$ such that we can bound the information complexity by

$$n(\varepsilon, d) \le C\varepsilon^{-p}d^q$$
 for all $d \in \mathbb{N}, \varepsilon \in (0, 1].$ (1.7)

If this last inequality holds with q = 0, i.e. if we have no dependence on the dimension at all, then the problem is called *strongly polynomially tractable*. In this case the smallest possible constant p in (1.7) is denoted by p^* . It is called the *exponent of strong polynomial tractability*.

If, in contrast, there do not exist constants C, p and q which fulfill (1.7) then the problem is said to be *polynomially intractable*.

Observe that (1.6) shows that, as long as the absolute error criterion is concerned, it is enough to consider $\varepsilon \in (0, \min \{\varepsilon_d^{\text{init}}, 1\}]$ instead of $\varepsilon \in (0, 1]$ in all the above definitions.

2. Properties and tools for special problem classes

This chapter deals with basic properties of certain classes of problems and algorithms. We state simple consequences obtained from fundamental assumptions on the operators under consideration. Furthermore, we present more or less classical tools used in the framework of information-based complexity to acquire tractability results in a quite general context.

In detail, we begin with a simple lower error bound in a very general setting which will be used on several occasions later on. In Section 2.2 we then show that for our purposes it is reasonable to concentrate mainly on compact problems and linear, non-adaptive algorithms. Moreover, there we derive a formula for the initial error of the problems we are interested in. Afterwards, in Section 2.3, we turn to the important class of problems defined between Hilbert spaces. We recall well-known tools such as the singular value decomposition, conclude optimal algorithms, and characterize several types of tractability of such problems. In Section 2.4 we restrict ourselves further and assume an additional tensor product structure which will play an important role throughout the rest of this paper. Finally we conclude this chapter with the discussion of so-called reproducing kernel Hilbert spaces.

The main references for the functional-analytic background needed in this part, as well as for the theory of s-numbers (or n-widths, respectively), are the monographs of Pinkus [Pin85] and Pietsch [Pie87, Pie07]. For a detailed discussion of applications to tractability questions we refer again to Novak and Woźniakowski [NW08, NW10, NW12] and to Mathé [Mat90].

2.1. Lower bounds on linear subspaces. For the purpose of this chapter it is enough to study the worst case setting. In addition, we will only focus on the case where all the problem elements lie in some centered ball of the respective source space. In this section we present quite a general method to obtain lower bounds on the nth minimal error with respect to a wide class of algorithms. In contrast to the rest of this paper (where we will restrict ourselves basically to linear and compact problems) we present a result that holds for any homogeneous operator S between linear normed spaces \mathcal{F} and \mathcal{G} over the field of real numbers. That is, we first only assume that $S(\alpha \cdot f) = \alpha \cdot S(f)$ for every $f \in \mathcal{F}$ and all $\alpha \in \mathbb{R}$

We start by proving the following (modified) assertion of Borsuk and Ulam for linear normed spaces:

LEMMA 2.1 (Borsuk–Ulam). Let V be a linear normed space over \mathbb{R} with $0 < \dim V = s$ $< \infty$ and let $N: V \to \mathbb{R}^n$ be a continuous mapping for some $0 \le n < s$. Then for all $r \ge 0$ there exists an element $f^* \in V$ with $||f^*| |V|| = r$ such that $N(f^*) = N(-f^*)$.

Proof. Obviously, the cases n=0, i.e. $N\equiv 0$, and r=0 are trivial. Hence, let $n\in\mathbb{N}$ and r>0. Since $\dim V=s$ we find an isomorphism $T\colon V\to\mathbb{R}^s$ such that T and T^{-1} are linear and bounded. Hence, for every r>0 the set $\Omega_r=T(\operatorname{int}(B_r(V)))$ is an open, bounded and symmetric subset of \mathbb{R}^s which contains zero. Moreover, the function $g=N\circ T^{-1}\colon \partial\Omega_r\to\mathbb{R}^n$ is continuous. From the theorem of Borsuk–Ulam (cf. Deimling [Dei85, Corollary 4.2]) we deduce the existence of some $\boldsymbol{x}^*\in\partial\Omega_r$ with $g(\boldsymbol{x}^*)=g(-\boldsymbol{x}^*)$. The claim now follows by taking $f^*=T^{-1}\boldsymbol{x}^*$.

With this result in hand, we can prove a generalization of Lemma 1 in [Wei12b].

PROPOSITION 2.2. Suppose S is a homogeneous operator between linear normed spaces \mathcal{F} and \mathcal{G} . Further assume that $V \subset \mathcal{F}$ is a linear subspace with dimension $s \in \mathbb{N}$ and that there exists a constant $a \geq 0$ such that

$$a||f| \mathcal{F}|| \le ||S(f)| \mathcal{G}|| \quad \text{for all } f \in V.$$
 (2.1)

Then for every $0 \le n < s$, any algorithm $A_n \in \mathcal{A}^{n,\text{cont}} \cup \mathcal{A}^{n,\text{adapt}}$, and all $r \ge 0$,

$$\Delta^{\text{wor}}(A_n; S \colon B_r(\mathcal{F}) \to \mathcal{G}) = \sup_{f \in B_r(\mathcal{F})} \|S(f) - A_n(f) \mid \mathcal{G}\| \ge ar.$$
 (2.2)

In particular, the nth minimal worst case error (within the unit ball $\mathcal{B}(\mathcal{F})$ of \mathcal{F}) satisfies $e^{\mathrm{wor}}(n; S \colon \mathcal{B}(\mathcal{F}) \to \mathcal{G}) \geq a$ for all n < s.

Proof. It is well-known that for $A_n = \varphi_n \circ N_n \in \mathcal{A}^{n,\text{cont}} \cup \mathcal{A}^{n,\text{adapt}}$ with n < s there exists $f^* \in V$ such that $N_n(f^*) = N_n(-f^*)$ and $||f^*|| \mathcal{F}|| = r$.

Without loss of generality let us again assume $n \in \mathbb{N}$ and r > 0 to avoid triviality. Then, for $A_n \in \mathcal{A}^{n,\mathrm{cont}}$, the existence of f^* is a simple consequence of Lemma 2.1 since in this case N_n is continuous by definition. On the other hand, if $A_n \in \mathcal{A}^{n,\mathrm{adapt}}$ then the proof can be obtained by arguments from linear algebra. We follow Werschulz and Woźniakowski [WW09, Theorem 3.1] and search for a non-zero $g \in V$ such that $N_n(g) = 0$, i.e.

$$L_1(g) = 0,$$

 $L_2(g; 0) = 0,$
 \vdots
 $L_n(g; 0, ..., 0) = 0.$ (2.3)

Since dim V=s every $g\in V$ can be represented uniquely as a linear combination $g=\sum_{m=1}^s c_m b_m$ of at most s linearly independent basis elements b_m of V. Due to the imposed linearity of $L_j(\cdot;0,\ldots,0), j=1,\ldots,n$, system (2.3) can be reformulated as a system of n homogeneous linear equations in the s>n unknowns $\mathbf{c}=(c_m)_{m=1}^s\in\mathbb{R}^s$. Consequently, it possesses a non-trivial solution $\mathbf{c}^*=(c_m^*)_{m=1}^s$, which implies the existence of some $g^*\in V\setminus\{0\}$ with $N_n(g^*)=0$. Since with L_j also N_n is linear, we can easily construct f^* out of g^* .

Anyway, every such f^* satisfies $A_n(f^*) = A_n(-f^*)$. Using the norm properties in the target space \mathcal{G} and the homogeneity of S we obtain (2.2):

$$\begin{split} \Delta^{\text{wor}}(A_n; S \colon B_r(\mathcal{F}) \to \mathcal{G}) &\geq \max \left\{ \| S(\pm f^*) - A_n(\pm f^*) \mid \mathcal{G} \| \right\} \\ &= \max \left\{ \| S(f^*) \pm A_n(f^*) \mid \mathcal{G} \| \right\} \\ &\geq \frac{1}{2} (\| S(f^*) + A_n(f^*) \mid \mathcal{G} \| + \| S(f^*) - A_n(f^*) \mid \mathcal{G} \|) \\ &\geq \frac{1}{2} \| 2 S(f^*) \mid \mathcal{G} \| \geq a \| f^* \mid \mathcal{F} \| = ar. \end{split}$$

The remaining implication for the *n*th minimal error finally follows from the case r=1 by taking the infimum over all $A_n \in \mathcal{A}^{n,\mathrm{cont}} \cup \mathcal{A}^{n,\mathrm{adapt}}$.

At this point we stress that the case $r \neq 1$ in (2.2) might be useful only if we deal with non-homogeneous (and thus non-linear) algorithms A_n . Otherwise we clearly have

$$\Delta^{\text{wor}}(A_n; S \colon B_r(\mathcal{F}) \to \mathcal{G}) = r\Delta^{\text{wor}}(A_n; S \colon \mathcal{B}(\mathcal{F}) \to \mathcal{G})$$

for all $r \geq 0$ provided that $S: \mathcal{F} \to \mathcal{G}$ is homogeneous. The importance of (2.2) for $r \neq 1$ will be made clear in Section 4.3.2 when we deal with embeddings $\mathcal{P} \hookrightarrow \mathcal{F}$. There we obtain a lower bound for the worst case error of S on $\mathcal{B}(\mathcal{F})$ out of a lower bound on $B_r(\mathcal{P})$ using $r = \|\mathrm{id} \mid \mathcal{L}(\mathcal{P}, \mathcal{F})\|^{-1}$.

2.2. Linearity and compactness. In what follows we will exclusively consider linear continuous problems $S = (S_d)_{d \in \mathbb{N}}$. That is, we assume that every solution operator S_d given by (1.2) is the restriction of a bounded linear mapping between some Banach spaces defined over the field of real numbers (⁶). If we assume that the set of problem elements $\widetilde{\mathcal{F}}_d$ is some ball $B_r(\mathcal{F}_d)$, r > 0, in the source space then conversely every bounded mapping S_d that acts linearly on this set (⁷) can be uniquely extended to a continuous linear operator \widetilde{S}_d on the whole space \mathcal{F}_d , i.e. $\widetilde{S}_d \in \mathcal{L}(\mathcal{F}_d, \mathcal{G}_d)$. From this point of view S_d and \widetilde{S}_d can be identified with each other and thus we use the symbol S_d for both.

At first glance the linearity assumption seems to be very restrictive. On the other hand, the two most important problems, namely approximation and integration, are indeed of this type. Moreover, the linear case is much better understood than the non-linear one; an overwhelming percentage of work on IBC was done in this setting. For the sake of completeness we also mention so-called quasilinear problems and refer to [WW07] and [NW12, Chapter 28].

Since we are interested in algorithms which are easy (and cheap) to implement, we pay special attention to the family of linear and non-adaptive algorithms $\mathcal{A}_d^{n,\mathrm{lin}}(\Lambda)$ (see (1.5)). It is well-known that this choice is reasonable for many classes of problems, since it can be shown that under mild assumptions optimal algorithms are indeed linear and non-adaptive. General assertions of this type can be found in Traub, Wasilkowski and

⁽⁶⁾ In fact, for most of the following results completeness is not needed. Many of them even remain valid (at least up to constants) using only quasi-norms or p-norms, but for simplicity we restrict ourselves to the case of Banach spaces. Finally, for the ease of notation, we only consider spaces over \mathbb{R} .

⁽⁷⁾ That means $S_d(\alpha \cdot f + \beta \cdot g)$ equals $\alpha \cdot S_d(f) + \beta \cdot S_d(g)$ for every convex combination $\alpha \cdot f + \beta \cdot g$ of elements $f, g \in \widetilde{\mathcal{F}}_d$.

Woźniakowski [TWW88], as well as in Novak and Woźniakowski [NW08, Section 4.2]. We do not present these results explicitly, since for the problems we are interested in, our assertions already imply such optimality statements.

Furthermore, we focus on information maps which are linear and continuous, i.e. $\Lambda \subseteq \Lambda^{\text{all}}$. Observe that then $A_{n,d} \in \mathcal{L}(\mathcal{F}_d, \mathcal{G}_d)$ and $\text{rank}(A_{n,d}) \leq n$. Moreover, for $\widetilde{\mathcal{F}}_d = \mathcal{B}(\mathcal{F}_d)$ we obtain

$$\Delta^{\text{wor}}(A_{n,d}; S_d) = \Delta^{\text{wor}}(A_{n,d}; S_d : \widetilde{\mathcal{F}}_d \to \mathcal{G}_d) = \|S_d - A_{n,d} \mid \mathcal{L}(\mathcal{F}_d, \mathcal{G}_d)\|.$$

It seems natural to ask when problems of this type are solvable at all. We say a problem $S = (S_d)_{d \in \mathbb{N}}$ is solvable if for any fixed $d \in \mathbb{N}$ there exists a sequence of algorithms $A_{n,d} \in \mathcal{A}_d^{n,\operatorname{lin}}(\Lambda^{\operatorname{all}})$ such that their worst case errors $\Delta^{\operatorname{wor}}(A_{n,d};S_d)$ tend to zero as $n \to \infty$. Hence, S_d needs to be an element of $\overline{\mathcal{FR}}(\mathcal{F}_d,\mathcal{G}_d)$, the closure of the finite rank operators in $\mathcal{L}(\mathcal{F}_d,\mathcal{G}_d)$, which is a subset of $\mathcal{K}(\mathcal{F}_d,\mathcal{G}_d)$. Therefore solvable problems are necessarily compact so that we can restrict ourselves in the following to $S_d \in \mathcal{K}(\mathcal{F}_d,\mathcal{G}_d)$. Due to the celebrated result of Enflo [Enf73], the converse is not true in this generality. Indeed, there are compact problems which are not solvable since there exist Banach spaces \mathcal{G}_d which do not have the so-called approximation property. However, the following (incomplete) list shows that in the cases we are interested in every compact problem is solvable:

PROPOSITION 2.3. Let $S = (S_d)_{d \in \mathbb{N}}$ with $S_d \in \mathcal{K}(\mathcal{F}_d, \mathcal{G}_d)$ for all $d \in \mathbb{N}$. Then S is solvable if for every $d \in \mathbb{N}$ one of the following conditions applies:

- the source space \mathcal{F}_d is a Hilbert space, or
- the target space \mathcal{G}_d is a Hilbert space, or
- the target space \mathcal{G}_d is $L_{\infty}(\mathcal{X}, \mathfrak{a}, \mu)$ for an arbitrary measure space $(\mathcal{X}, \mathfrak{a}, \mu)$.

Proof. Let $d \in \mathbb{N}$. Given all the above restrictions we note that if we consider the class $\mathcal{A}_d^{n,\text{lin}}(\Lambda^{\text{all}})$ then the numbers $e^{\text{wor}}(n,d;S_d), n \in \mathbb{N}_0$, per definition equal the linear n-widths (or approximation numbers) $\delta_n(S_d)$ as defined in [Pin85, Definition 7.3]. Up to an index shift these numbers form an s-scale (8) in the sense of Pietsch [Pie07, Section 6.2]. Other important s-scales are the Gelfand numbers $c_n(S_d)$ and the Kolmogorov numbers $d_n(S_d)$. Without going into details we mention that for any compact operator $S_d \in \mathcal{K}(\mathcal{F}_d, \mathcal{G}_d)$ both these sequences tend to zero as $n \to \infty$ (see [Pin85, Propositions 7.4 and 7.1]). Hence, to prove solvability it suffices to show that $\delta_n(S_d) \leq \max\{c_n(S_d), d_n(S_d)\}$ for all $n \in \mathbb{N}$. Indeed, if \mathcal{F}_d is a Hilbert space then $\delta_n(S_d) = c_n(S_d)$. Furthermore, $\delta_n(S_d) = d_n(S_d)$ if \mathcal{G}_d is a Hilbert space (see, e.g., [Pin85, p. 33]). Finally Proposition 8.13 in [Pin85] shows that the second last equality remains valid if the target space \mathcal{G}_d enjoys the so-called (metric) extension property. It is known that in particular $L_\infty(\mathcal{X},\mathfrak{a},\mu)$ has this property (see, e.g., König [Kön86, 1.c.2]).

We stress that in Proposition 2.3 we do not need to assume that the Hilbert spaces are separable.

Let us conclude this section with a proposition which shows that the zero algorithm $A_0 \equiv 0$ is the optimal choice among all approximations to a given operator $S \in \mathcal{L}(\mathcal{F}, \mathcal{G})$

⁽⁸⁾ Note that for historical reasons there is some notational confusion concerning s-numbers and n-widths. See, e.g., [Pie07, p. 336] for details.

that do not use any information on the input $f \in \mathcal{F}$. Here \mathcal{F} and \mathcal{G} can be arbitrary normed spaces.

PROPOSITION 2.4. For $S \in \mathcal{L}(\mathcal{F}, \mathcal{G})$ and $\mathcal{A}_0 = 0 \in \mathcal{L}(\mathcal{F}, \mathcal{G})$ we have

$$e^{\text{wor}}(0; S \colon \mathcal{B}(\mathcal{F}) \to \mathcal{G}) = \Delta^{\text{wor}}(A_0; S \colon \mathcal{B}(\mathcal{F}) \to \mathcal{G}) = ||S| \mathcal{L}(\mathcal{F}, \mathcal{G})||.$$

Consequently, the zero algorithm is optimal for S within the class $\mathcal{A}^{0,\mathrm{cont}} \cup \mathcal{A}^{0,\mathrm{adapt}}$ and the initial worst case error $\varepsilon^{\mathrm{init,wor}}$ of S is given by its operator norm.

Proof. Obviously the second equality is true by the definition of Δ^{wor} . Moreover, the linear algorithm $A_0 \equiv 0$ is included in every class of algorithms we defined in Section 1.3. This implies in particular $e^{\text{wor}}(0; S) \leq \Delta^{\text{wor}}(A_0) = ||S| \mathcal{L}(\mathcal{F}, \mathcal{G})||$.

To show the converse inequality, recall that every algorithm A that does not use any information on the input necessarily takes the form $A(f) \equiv g$ for some $g \in \mathcal{G}$. A calculation similar to that in the proof of Proposition 2.2 yields

$$||S(f)||\mathcal{G}|| \le \max\{||S(f) - g||\mathcal{G}||, ||S(-f) - g||\mathcal{G}||\}$$

for every $f \in \mathcal{F}$. Taking the supremum over $f \in \mathcal{B}(\mathcal{F})$ now shows that

$$||S \mid \mathcal{L}(\mathcal{F}, \mathcal{G})|| \leq \Delta^{\text{wor}}(A),$$

which implies the desired result since A was chosen arbitrarily.

We note in passing that the last step in the above proof crucially depends on the fact that the unit ball $\widetilde{\mathcal{F}} = \mathcal{B}(\mathcal{F})$ is symmetric in the sense that $f \in \widetilde{\mathcal{F}}$ implies $-f \in \widetilde{\mathcal{F}}$.

- **2.3.** General Hilbert space problems. In this section we describe the *singular value decomposition* (SVD), which turns out to be the main tool when dealing with problems where both the source and the target spaces are Hilbert spaces. We prove well-known formulas for optimal linear algorithms using continuous linear functionals and calculate their worst case errors. Afterwards, we use the assertions obtained to give characterizations of (strong) polynomial tractability for these problems.
- **2.3.1. Singular value decomposition.** Given any compact operator $T \in \mathcal{K}(\mathcal{F}, \mathcal{G})$ acting between real Hilbert spaces \mathcal{F} and \mathcal{G} we define its *adjoint operator* $T^{\dagger} \colon \mathcal{G} \to \mathcal{F}$ in the usual way by

$$\langle Tf, g \rangle_{\mathcal{G}} = \langle f, T^{\dagger}g \rangle_{\mathcal{F}} \quad \text{ for all } f \in \mathcal{F}, g \in \mathcal{G}.$$
 (2.4)

Of course, T^{\dagger} is always unique and well-defined. For details we refer the reader to Yosida [Yos80, VII.2]. If $\mathcal{F} = \mathcal{G}$ and $T^{\dagger} = T$, then we say that T is *self-adjoint*. From Schauder's theorem we know that $T^{\dagger} \in \mathcal{K}(\mathcal{G}, \mathcal{F})$ if and only if $T \in \mathcal{K}(\mathcal{F}, \mathcal{G})$ (see e.g. [Pin85, p. 31]). Hence, it is easily seen that also

$$W = T^{\dagger}T \colon \mathcal{F} \to \mathcal{F}$$

defines a compact operator. Moreover, W is obviously self-adjoint and positive, i.e. $\langle Wf, f \rangle_{\mathcal{F}} \geq 0$ for every $f \in \mathcal{F}$. It is well-known that then all the eigenvalues $\lambda_m = \lambda_m(W)$ of W are necessarily real and non-negative. Following Pinkus [Pin85, p. 64] we denote the sum of the algebraic multiplicities of the non-zero eigenvalues of W by v = v(W). Note that Riesz–Schauder theory proves that there are at most countably many non-zero

eigenvalues. They are uniformly bounded, each has a finite multiplicity and there are no accumulation points but (possibly) zero (see, e.g., [Yos80, X.5, Theorem 2]). Observe further that in any case $v \leq \dim \mathcal{F} \in \mathbb{N} \cup \{\infty\}$. Let us write these eigenvalues in non-increasing ordering subscripted by indices from the set $\mathcal{M} = \{m \in \mathbb{N} \mid m < v + 1\}$:

$$\lambda_1 \ge \dots \ge \lambda_m \ge \dots > 0. \tag{2.5}$$

Without loss of generality we will always assume the existence of at least one non-trivial eigenvalue, i.e. we exclude the operator $T \equiv 0$, which ensures that $\mathcal{M} \neq \emptyset$. We denote the corresponding (mutually orthonormal) eigenvectors of W by ϕ_m , $m \in \mathcal{M}$, and refer to $\{(\lambda_m, \phi_m) \mid m \in \mathcal{M}\}$ as the set of non-trivial eigenpairs of W. Consequently, for $i, j \in \mathcal{M}$, by (2.4) we have

$$\langle T\phi_i, T\phi_j \rangle_{\mathcal{G}} = \langle \phi_i, T^{\dagger}T\phi_j \rangle_{\mathcal{F}} = \langle \phi_i, W\phi_j \rangle_{\mathcal{F}} = \langle \phi_i, \lambda_j \phi_j \rangle_{\mathcal{F}} = \delta_{i,j} \cdot \lambda_j. \tag{2.6}$$

If we extend the possibly finite eigenvalue sequence $(\lambda_m)_{m=1}^v$ by taking $\lambda_m = 0$ for all m > v then clearly $\lambda = (\lambda_m)_{m \in \mathbb{N}}$ forms a null sequence. Again following Pinkus, we call the square root $\sigma = \sigma(T)$ of $\lambda = \lambda(W)$,

$$\sigma_m = \sqrt{\lambda_m}, \quad m \in \mathbb{N},$$

the sequence of $singular\ values$ of T. The importance of this bunch of definitions comes from the following assertion.

Theorem 2.5 (Singular value decomposition). Let \mathcal{F} and \mathcal{G} be arbitrary Hilbert spaces and $T \in \mathcal{K}(\mathcal{F}, \mathcal{G})$. Then, with the above notations,

$$T = \sum_{m=1}^{v} \langle \cdot, \phi_m \rangle_{\mathcal{F}} T \phi_m. \tag{2.7}$$

Proof. A detailed proof can be found in the monograph of König [Kön86, 1.b.3]. It is mainly based on the so-called *polar decomposition* of continuous linear operators and Riesz–Schauder theory. Actually, the proof deals with complex Hilbert spaces but it literally transfers to the real case. Moreover, only the existence of an orthonormal sequence $(\psi_m)_{m=1}^v$ in \mathcal{G} is shown such that pointwise,

$$Tf = \sum_{m=1}^{v} \sigma_m \langle f, \phi_m \rangle_{\mathcal{F}} \psi_m, \quad f \in \mathcal{F}.$$

However, setting $f = \phi_k$ for $k \in \mathcal{M}$ together with the mutual orthonormality of $(\phi_m)_{m=1}^v$ immediately implies $\sigma_k \psi_k = T \phi_k$ for any k. The claimed identity in $\mathcal{K}(\mathcal{F}, \mathcal{G})$ finally follows from Bessel's inequality.

REMARK 2.6. Note that again the Hilbert spaces \mathcal{F} and \mathcal{G} do not need to be separable. Nevertheless the image of \mathcal{F} under T is indeed separable, because it is spanned by at most countably many elements $T\phi_m \in \mathcal{G}$. Since the elements of the set $\Phi = \{\phi_m \in \mathcal{F} \mid m \in \mathcal{M}\}$ are mutually orthonormal we can extend Φ to an orthonormal basis (ONB) E of \mathcal{F} . Then (2.7) shows that $\ker T = \Phi^{\perp}$. Remember that we are only interested in the approximation of the image of T. Hence, we can without loss of generality assume that $E = \Phi$. In other words, even though \mathcal{F} may be non-separable, in general we can restrict ourselves

to the separable case in what follows. We only need to replace \mathcal{F} by $\overline{\Phi}$, the closure of the orthonormal eigenelements of $W = T^{\dagger}T$ under $\langle \cdot, \cdot \rangle_{\mathcal{F}}$. \square

2.3.2. Optimal algorithm. Observe that by (2.7) we obtained a representation of any operator $T \in \mathcal{K}(\mathcal{F}, \mathcal{G})$ as the limit of related finite rank operators. Therefore we are able to construct nth optimal linear algorithms which only use information from Λ^{all} . This is stated in the following corollary which can be found (slightly modified) as Corollary 4.12 in [NW08].

COROLLARY 2.7. For $d \in \mathbb{N}$ assume \mathcal{F}_d and \mathcal{G}_d are arbitrary Hilbert spaces. Further let $S = (S_d)_{d \in \mathbb{N}}$ denote a compact problem acting between these spaces, i.e. $S_d \in \mathcal{K}(\mathcal{F}_d, \mathcal{G}_d)$ for every d. Then for all $d \in \mathbb{N}$ and $n \in \mathbb{N}_0$ the algorithm $A_{n,d}^* \in \mathcal{A}_d^{n,\mathrm{lin}}(\Lambda^{\mathrm{all}})$ given by

$$A_{n,d}^* \colon \mathcal{F}_d \to \mathcal{G}_d, \quad f \mapsto A_{n,d}^*(f) = \sum_{m=1}^{\min\{n,v(W_d)\}} \langle f, \phi_{d,m} \rangle_{\mathcal{F}_d} S_d \phi_{d,m},$$

for S_d is optimal in the class $\mathcal{A}_d^{n,\mathrm{cont}} \cup \mathcal{A}_d^{n,\mathrm{adapt}}$ and we have

$$e^{\text{wor}}(n, d; S_d) = \Delta^{\text{wor}}(A_{n,d}^*; S_d) = \sigma_{d,n+1} = \sqrt{\lambda_{d,n+1}}.$$
 (2.8)

Here for $d \in \mathbb{N}$ the singular values $(\sigma_{d,m})_{m \in \mathbb{N}}$, as well as the eigenvectors $(\phi_{d,m})_{m=1}^{v(W_d)}$, are constructed from $W_d = S_d^{\dagger} S_d$ as explained above.

Proof. Recall that $\Delta^{\text{wor}}(A_{n,d}^*; S_d)$ equals $||S_d - A_{n,d}^*| | \mathcal{L}(\mathcal{F}_d, \mathcal{G}_d)||$ for any fixed $d \in \mathbb{N}$ and $n \in \mathbb{N}_0$. Without loss of generality we can assume $n < v = v(W_d)$ since otherwise $A_{n,d}^* = S_d$ due to (2.7). This would imply (2.8) because $\sigma_{d,m} = 0$ for all m > v.

Let $M \in \mathbb{N}$ with $n+1 \leq M \leq v$ and fix $f \in \mathcal{B}(\mathcal{F}_d)$. Then, by (2.6), the non-increasing ordering of $(\lambda_m)_{m=1}^v$, and Bessel's inequality,

$$\left\| \sum_{m=n+1}^{M} \langle f, \phi_{d,m} \rangle_{\mathcal{F}_d} S_d \phi_{d,m} \mid \mathcal{G}_d \right\|^2 = \sum_{m=n+1}^{M} \langle f, \phi_{d,m} \rangle_{\mathcal{F}_d}^2 \lambda_{d,m} \leq \lambda_{d,n+1} \sum_{m=1}^{v} \langle f, \phi_{d,m} \rangle_{\mathcal{F}_d}^2$$
$$\leq \lambda_{d,n+1} \|f \mid \mathcal{F}_d \|^2 \leq \sigma_{d,n+1}^2.$$

The particular choice $f = \phi_{d,n+1}$ shows that the estimates are sharp. Anyway, we obtain $||S_d - A_{n,d}^*| \mathcal{L}(\mathcal{F}_d, \mathcal{G}_d)|| \leq \sigma_{d,n+1}$, which proves

$$e^{\text{wor}}(n, d; S_d) \le \Delta^{\text{wor}}(A_{n,d}^*; S_d) \le \sigma_{d,n+1}.$$

To show the converse, i.e. $e^{\text{wor}}(n,d;S_d) \geq \sigma_{d,n+1}$ for $n \in \mathbb{N}_0$ and $d \in \mathbb{N}$, we use Parseval's identity on $V = \text{span}\{\phi_{d,m} \mid m \leq n+1\} \subset \mathcal{F}_d$ together with the linearity of S_d to obtain

$$||S_d f | \mathcal{G}_d||^2 = \sum_{m=1}^{n+1} \langle f, \phi_{d,m} \rangle_{\mathcal{F}_d}^2 \lambda_{d,m} \ge \lambda_{d,n+1} \sum_{m=1}^{n+1} \langle f, \phi_{d,m} \rangle_{\mathcal{F}_d}^2 = \sigma_{d,n+1}^2 ||f| |\mathcal{F}_d||^2$$

for all $f \in V$. The claim now follows from the application of Proposition 2.2 with $a = \sigma_{d,n+1}$. Moreover, Proposition 2.2 also shows that we cannot reduce the error by taking algorithms $A_{n,d} \in (\mathcal{A}_d^{n,\mathrm{cont}} \cup \mathcal{A}_d^{n,\mathrm{adapt}}) \setminus \mathcal{A}_d^{n,\mathrm{lin}}$.

Note that (2.8) together with Proposition 2.4 implies in particular that for $d \in \mathbb{N}$ the initial worst case error of S_d is

$$\varepsilon_d^{\text{init,wor}} = ||S_d | \mathcal{L}(\mathcal{F}_d, \mathcal{G}_d)|| = \sigma_{d,1} = \sqrt{\lambda_{d,1}}.$$

2.3.3. Polynomial tractability. As an immediate consequence of (2.8) we can calculate the information complexity of Hilbert space problems in the worst case setting (with respect to the class $\mathcal{A}_d^{n,\mathrm{lin}}(\Lambda^{\mathrm{all}})$) for every $d \in \mathbb{N}$ and $\varepsilon > 0$ by

$$n_{\text{abs}}^{\text{wor}}(\varepsilon, d) = \min \left\{ n \in \mathbb{N}_0 \mid \sigma_{d, n+1} \le \varepsilon \right\} = \# \left\{ n \in \mathbb{N} \mid \lambda_{d, n} > \varepsilon^2 \right\}$$
 (2.9)

for the absolute error criterion, and by

$$n_{\text{norm}}^{\text{wor}}(\varepsilon, d) = \#\{n \in \mathbb{N} \mid \lambda_{d,n}/\lambda_{d,1} > \varepsilon^2\}$$
 (2.10)

for the normalized error criterion. This observation leads to the following refinement of Theorem 5.1 in Novak and Woźniakowski [NW08], which can also be found in [Wei11]. It gives necessary and sufficient conditions for (strong) polynomial tractability in terms of summability properties of the sequences $(\lambda_{d,m})_{m\in\mathbb{N}}$.

Theorem 2.8. Assume S is a problem as in Corollary 2.7 and consider the absolute error criterion in the worst case setting.

• If S is polynomially tractable with the constants C, p > 0 and $q \ge 0$ then for all $\tau > p/2$ we have

$$C_{\tau} = \sup_{d \in \mathbb{N}} \frac{1}{d^r} \left(\sum_{i=f(d)}^{\infty} \lambda_{d,i}^{\tau} \right)^{1/\tau} < \infty, \tag{2.11}$$

where r=2q/p and $f:\mathbb{N}\to\mathbb{N}$ with $f(d)=\lceil (1+C)\,d^q\rceil$. In this case $C_\tau\leq C^{2/p}\cdot \zeta(2\tau/p)^{1/\tau}$.

• If (2.11) is satisfied for some parameters $r \geq 0$, $\tau > 0$ and a function $f \colon \mathbb{N} \to \mathbb{N}$ such that $f(d) = \lceil C(\min\left\{\varepsilon_d^{\text{init}}, 1\right\})^{-p} d^q \rceil$, where C > 0 and $p, q \geq 0$, then the problem S is polynomially tractable. In detail, we have the bound $n_{\text{abs}}^{\text{wor}}(\varepsilon, d) \leq (C + C_{\tau}^{\tau}) \varepsilon^{-\max\{p, 2\tau\}} d^{\max\{q, r\tau\}}$ for any $\varepsilon \in (0, 1]$ and every $d \in \mathbb{N}$.

Proof. If the problem is polynomially tractable then there exist constants C, p > 0 and $q \ge 0$ such that for all $d \in \mathbb{N}$ and $\varepsilon \in (0, 1]$,

$$n(\varepsilon, d) = n_{\text{abs}}^{\text{wor}}(\varepsilon, d) \le C \varepsilon^{-p} d^q.$$

Formula (2.9) and the non-increasing ordering of $(\lambda_{d,i})_{i\in\mathbb{N}}$ therefore imply

$$\lambda_{d,\lfloor C\varepsilon^{-p}d^q\rfloor+1} \le \lambda_{d,n(\varepsilon,d)+1} \le \varepsilon^2, \quad \varepsilon \in (0,1].$$

If we set $i = \lfloor C\varepsilon^{-p}d^q \rfloor + 1$ and vary ε in (0,1] then i takes the values $\lfloor C d^q \rfloor + 1$, $\lfloor C d^q \rfloor + 2$, and so forth. On the other hand, we have $i \leq C\varepsilon^{-p}d^q + 1$, which is equivalent to $\varepsilon^2 \leq (Cd^q/(i-1))^{2/p}$ if $i \geq 2$. For all $i \geq f(d) = \lceil (1+C)d^q \rceil$ we indeed have $i \geq 2$, and consequently

$$\lambda_{d,i} \le \lambda_{d,n(\varepsilon,d)+1} \le \varepsilon^2 \le \left(\frac{Cd^q}{i-1}\right)^{2/p}.$$

Choosing $\tau > p/2 > 0$ we conclude that

$$\sum_{i=f(d)}^{\infty} \lambda_{d,i}^{\tau} \leq \sum_{i=f(d)}^{\infty} \left(\frac{Cd^q}{i-1}\right)^{2\tau/p} = (Cd^q)^{2\tau/p} \sum_{i=f(d)-1}^{\infty} \frac{1}{i^{2\tau/p}} \leq \left(C^{2/p}d^{2q/p}\right)^{\tau} \zeta\left(\frac{2\tau}{p}\right)^{2\tau/p}$$

for every $d \in \mathbb{N}$. In other words, we have shown (2.11) with r = 2q/p, as well as the estimate on C_{τ} .

Conversely, assume that for some $r \geq 0$ and $\tau > 0$ estimate (2.11) holds with

$$f(d) = \left[C(\min\left\{ \varepsilon_d^{\text{init}}, 1 \right\})^{-p} d^q \right], \text{ where } C > 0 \text{ and } p, q \ge 0.$$

That is, we assume $0 < C_{\tau} < \infty$. For $n \geq f(d)$ the ordering of $(\lambda_{d,i})_{i \in \mathbb{N}}$ implies $\sum_{i=f(d)}^{n} \lambda_{d,i}^{\tau} \geq \lambda_{d,n}^{\tau} (n-f(d)+1)$. Hence, for every $d \in \mathbb{N}$ and $n \geq f(d)$,

$$\lambda_{d,n}(n-f(d)+1)^{1/\tau} \le \left(\sum_{i=f(d)}^{n} \lambda_{d,i}^{\tau}\right)^{1/\tau} \le \left(\sum_{i=f(d)}^{\infty} \lambda_{d,i}^{\tau}\right)^{1/\tau} \le C_{\tau} d^{r},$$

or respectively, $\lambda_{d,n+1} \leq C_\tau d^r ((n+1) - f(d) + 1)^{-1/\tau}$, for all $n \geq f(d) - 1$. Note that for $\varepsilon \in (0, \min\{\varepsilon_d^{\text{init}}, 1\}]$ we have $C_\tau d^r ((n+1) - f(d) + 1)^{-1/\tau} \leq \varepsilon^2$ if and only if

$$n \ge n^* = \left\lceil \left(\frac{C_\tau d^r}{\varepsilon^2} \right)^\tau \right\rceil + f(d) - 2.$$

In particular, $\lambda_{d,n+1} \leq \varepsilon^2$ at least for $n \geq \max\{n^*, f(d) - 1\}$. In other words, for every $d \in \mathbb{N}$ and all $\varepsilon \in (0, \min\{\varepsilon_d^{\text{init}}, 1\}]$,

$$n_{\text{abs}}^{\text{wor}}(\varepsilon, d) \le \max\{n^*, f(d) - 1\} \le f(d) - 1 + \left(\frac{C_{\tau}}{\varepsilon^2}\right)^{\tau}$$

$$\le C\left(\min\left\{\varepsilon_d^{\text{init}}, 1\right\}\right)^{-p} d^q + C_{\tau}^{\tau} \varepsilon^{-2\tau} d^{\tau\tau} \le \left(C + C_{\tau}^{\tau}\right) \varepsilon^{-\max\{p, 2\tau\}} d^{\max\{q, r\tau\}}.$$

Thus, the problem is polynomially tractable since $n_{\rm abs}^{\rm wor}(\varepsilon,d)=0$ for $\varepsilon\geq \varepsilon_d^{\rm init}$.

Let us add some comments on this result. Theorem 2.8 clearly provides a characterization of (strong) polynomial tractability. In comparison to Theorem 5.1 in [NW08] our result has the essential advantage that the estimates incorporate the initial error $\varepsilon_d^{\rm init}$. Hence if $\varepsilon_d^{\rm init}$ is sufficiently small then we can deduce polynomial tractability while ignoring a larger set of eigenvalues in the summation (2.11).

Observe that the first statement does not cover any assertion about the initial error itself, since $f(d) \geq 2$. Thus it might happen that we have (strong) polynomial tractability with respect to the absolute error criterion, though the largest eigenvalue $\lambda_{d,1} = (\varepsilon_d^{\text{init}})^2$ tends to infinity faster than any polynomial. To give an example, for $d \in \mathbb{N}$ we consider the sequences $(\lambda_{d,m})_{m \in \mathbb{N}}$ defined by

$$\lambda_{d,1} = e^{2d}$$
 and $\lambda_{d,m} = \frac{1}{m}$ for $m \ge 2$.

Here, obviously, the initial error grows to infinity exponentially fast, but nevertheless the second point of Theorem 2.8 shows that S is strongly polynomially tractable since (2.11) holds with r = p = q = 0, and $C = \tau = 2$.

Next we present an analogue of Theorem 2.8 for the normalized error criterion. Again a slightly modified statement can be found in [NW08, Theorem 5.2].

THEOREM 2.9. Assume S is as in Corollary 2.7 and consider the normalized error criterion in the worst case setting.

• If S is polynomially tractable with the constants C, p > 0 and $q \ge 0$ then for all $\tau > p/2$ we have

$$C_{\tau} = \sup_{d \in \mathbb{N}} \frac{1}{d^r} \left(\sum_{i=f(d)}^{\infty} \left(\frac{\lambda_{d,i}}{\lambda_{d,1}} \right)^{\tau} \right)^{1/\tau} < \infty, \tag{2.12}$$

where r = 2q/p and $f: \mathbb{N} \to \mathbb{N}$ with $f(d) \equiv 1$. In this case the bound $C_{\tau} \leq 2^{1/\tau} (1+C)^{2/p} \zeta(2\tau/p)^{1/\tau}$ holds for any such τ .

• If (2.11) is satisfied for some parameters $r \geq 0$, $\tau > 0$ and a function $f \colon \mathbb{N} \to \mathbb{N}$ such that $f(d) = \lceil Cd^q \rceil$, where C > 0 and $q \geq 0$, then the problem S is polynomially tractable. If so, then $n_{\text{norm}}^{\text{wor}}(\varepsilon, d) \leq (C + C_{\tau}^{\tau}) \varepsilon^{-2\tau} d^{\max\{q, r\tau\}}$ for any $\varepsilon \in (0, 1]$ and every $d \in \mathbb{N}$.

Proof. By the strong relation between the absolute and the normalized error criteria, i.e. $n_{\text{norm}}^{\text{wor}}(\varepsilon,d) = n_{\text{abs}}^{\text{wor}}(\varepsilon\sqrt{\lambda_{d,1}},d)$ for $\varepsilon \in (0,1]$ and $d \in \mathbb{N}$, we note that Theorem 2.9 can be shown using essentially the same arguments as in the proof of Theorem 2.8. Indeed, if we replace $\lambda_{d,i}$ by $\lambda_{d,i}/\lambda_{d,1}$ for $i \in \mathbb{N}$ we obtain a scaled problem T with initial error $\varepsilon_d^{\text{init}} = 1$. Now the information complexity of T (with respect to the absolute error criterion) equals the information complexity of S with respect to normalized errors (9). Following the lines of the proof of Theorem 2.8 this shows the second point of Theorem 2.9, where we set p = 0. Moreover, for any $\tau > p/2$ and $d \in \mathbb{N}$ we conclude that

$$\sum_{i=\lceil (1+C)d^q \rceil}^{\infty} \left(\frac{\lambda_{d,i}}{\lambda_{d,1}}\right)^{\tau} \leq (C^{2/p} d^{2q/p})^{\tau} \zeta\left(\frac{2\tau}{p}\right) \leq (1+C)^{2\tau/p} \zeta\left(\frac{2\tau}{p}\right) d^{q\cdot 2\tau/p},$$

provided that S is polynomially tractable with constants C, p > 0 and $q \ge 0$. Furthermore, for any $d \in \mathbb{N}$ we have

$$\sum_{i=1}^{\lceil (1+C)d^q \rceil - 1} \left(\frac{\lambda_{d,i}}{\lambda_{d,1}} \right)^{\tau} \le \lceil (1+C)d^q \rceil - 1 \le (1+C) d^q \le (1+C)^{2\tau/p} \zeta \left(\frac{2\tau}{p} \right) d^{q \cdot 2\tau/p}$$

since $\lambda_{d,i} \leq \lambda_{d,1}$, $2\tau/p > 1$, and $\zeta(2\tau/p) > 1$. Consequently, setting r = 2q/p and combining both the previous estimates leads to

$$\frac{1}{d^r} \left(\sum_{i=1}^{\infty} \left(\frac{\lambda_{d,i}}{\lambda_{d,1}} \right)^{\tau} \right)^{1/\tau} \le 2^{1/\tau} \left(1 + C \right)^{2/p} \zeta \left(\frac{2\tau}{p} \right)^{1/\tau} \quad \text{for } d \in \mathbb{N},$$

which shows (2.12), as well as the claimed bound on C_{τ} .

Obviously Theorem 2.9 again provides a characterization of (strong) polynomial tractability of a given compact Hilbert space problem $S = (S_d)_{d \in \mathbb{N}}$ in terms of summability properties of the eigenvalue sequence $(\lambda_{d,i})_{i \in \mathbb{N}}$ of $W_d = S_d^{\dagger} S_d$.

⁽⁹⁾ For details we refer to the proof of Theorem 2.12.

- **2.4. Tensor product problems.** In the previous section we investigated tractability properties of compact Hilbert space problems $S = (S_d)_{d \in \mathbb{N}}$ without assuming any relation between subsequent problem operators S_d . Next we want to consider problems S where every S_d is generated out of one single (univariate) operator S_1 via a d-fold tensor product construction.
- **2.4.1. Definition and simple properties.** We need to recall the concept of tensor product Hilbert spaces. To this end, we use the approach of Kadison and Ringrose [KR83, Section 2.6]. For a comprehensive introduction to more general tensor products in functional analysis we refer to the first chapter of Light and Cheney [LC85] and to Section 1.3 in Hansen [Han10].

Without going into detail, we note that given a finite number of arbitrary Hilbert spaces $H^{(k)}$ with inner products $\langle \cdot, \cdot \rangle_{H^{(k)}}$, $k = 1, \ldots, d$, the tensor product space

$$H_d = \bigotimes_{k=1}^d H^{(k)} = H^{(1)} \otimes \cdots \otimes H^{(d)}$$

can be identified (10) with the closure of the algebraic tensor product $H_{d,0}$, with respect to a (reasonable cross) norm which is induced by a certain inner product $\langle \cdot, \cdot \rangle_{H_{d,0}}$. Keep in mind that the algebraic tensor product is defined as the quotient of the *free vector space*, i.e. the set of all finite linear combinations of formal objects $f = \bigotimes_{k=1}^{d} f_k$ with $f_k \in H^{(k)}$, which we call *simple* (or *pure*) *tensors*, by a suitable linear subspace (11). Moreover, the inner product on the algebraic tensor product $H_{d,0}$ just mentioned is defined by

$$\left\langle \bigotimes_{k=1}^{d} f_k, \bigotimes_{k=1}^{d} g_k \right\rangle_{H_{d,0}} = \prod_{k=1}^{d} \left\langle f_k, g_k \right\rangle_{H^{(k)}} \quad \text{for } f_k, g_k \in H^{(k)}.$$

By means of continuous (multi)linear extension this functional uniquely determines the inner product $\langle \cdot, \cdot \rangle_{H_d}$ on H_d . As usual we denote the corresponding norm by $\|\cdot\|_{H_d}$.

Due to the tensor product structure, many useful properties such as completeness and separability of the underlying spaces $H^{(k)}$ are transferred directly to H_d provided that all the $H^{(k)}$ share them. In particular, it is well-known how to construct an orthonormal basis (ONB) of the tensor product space given an ONB

$$E^{(k)} = \{ e_i^{(k)} \in H^{(k)} \mid i \in \mathcal{I}^{(k)} \}$$

in each $H^{(k)}$, k = 1, ..., d. Here every $\mathcal{I}^{(k)}$ denotes a (possibly uncountable) abstract index set. Then the set of all d-fold simple tensors given by

$$E_d = \left\{ e_{d,\boldsymbol{j}} = \bigotimes_{k=1}^d e_{j_k}^{(k)} \mid \boldsymbol{j} = (j_1, \dots, j_d) \in \mathcal{I}_d = \bigotimes_{k=1}^d \mathcal{I}^{(k)} \right\}$$

is the desired ONB in H_d (see [KR83, Theorem 2.6.4]).

For the applications we have in mind we will focus on the special case where all the building blocks $H^{(k)}$, k = 1, ..., d, of H_d coincide. In what follows we therefore assume

 $[\]binom{10}{10}$ Note that this association is unique up to some isometric isomorphism.

⁽¹¹⁾ To abbreviate the notation we do not distinguish between simple tensors and their equivalence classes in what follows.

 $H^{(k)} \equiv H_1$ for some Hilbert space H_1 . The respective ONB of H_1 will be denoted by $E_1 = \{e_i \in H_1 \mid i \in \mathcal{I}_1\}$. Then the above formula for E_d simplifies to

$$E_d = \left\{ e_{d,j} = \bigotimes_{k=1}^d e_{j_k} \mid j = (j_1, \dots, j_d) \in \mathcal{I}_d = (\mathcal{I}_1)^d \right\}.$$
 (2.13)

We are ready to introduce the tensor product problem operators S_d , $d \geq 1$, we are interested in. Let $S_1 \colon \mathcal{F}_1 \to \mathcal{G}_1$ be a compact linear operator between arbitrary Hilbert spaces \mathcal{F}_1 and \mathcal{G}_1 . For $d \geq 2$ we assume $\mathcal{F}_d = H_d$ is the d-fold tensor product space of $H^{(k)} = H_1 = \mathcal{F}_1$, $k = 1, \ldots, d$, as explained above. Analogously, we construct the space $\mathcal{G}_d = \bigotimes_{k=1}^d \mathcal{G}_1$ out of d copies of \mathcal{G}_1 . Now Proposition 2.6.12 of [KR83] implies that there exists a uniquely defined linear operator $S_d = \bigotimes_{k=1}^d S_1 \colon \mathcal{F}_d \to \mathcal{G}_d$ such that

$$S_d\left(\bigotimes_{k=1}^d f_k\right) = \bigotimes_{k=1}^d S_1 f_k, \quad f_k \in \mathcal{F}_1,$$

and we have $||S_d| | \mathcal{L}(\mathcal{F}_d, \mathcal{G}_d)|| = ||S_1| | \mathcal{L}(\mathcal{F}_1, \mathcal{G}_1)||^d < \infty$ for any fixed $d \in \mathbb{N}$. In detail, we define the bounded linear operator $\widetilde{S}_d \colon E_d \to \mathcal{G}_d$ such that for all $\mathbf{j} \in \mathcal{I}_d$ we have $\widetilde{S}_d(e_{d,\mathbf{j}}) = \widetilde{S}_d(\bigotimes_{k=1}^d e_{j_k}) = \bigotimes_{k=1}^d S_1(e_{j_k}) \in \mathcal{G}_d$. Then S_d is assumed to be the uniquely defined continuous linear extension of \widetilde{S}_d from E_d to \mathcal{F}_d . By the compactness of S_1 it is easy to check that the problem operator S_d is not only bounded but even compact. Moreover, a linear extension argument shows that the adjoint operator S_d^{\dagger} is the d-fold tensor product of S_1^{\dagger} , i.e. $S_d^{\dagger} = \bigotimes_{k=1}^d S_1^{\dagger}$, and hence

$$W_d = S_d^{\dagger} S_d = \left(\bigotimes_{k=1}^d S_1^{\dagger} \right) \left(\bigotimes_{k=1}^d S_1 \right) = \bigotimes_{k=1}^d S_1^{\dagger} S_1 = \bigotimes_{k=1}^d W_1$$
 (2.14)

(cf. [KR83, p. 146]).

2.4.2. Eigenpairs and the optimal algorithm. From Section 2.3 we know that for d in \mathbb{N} the optimal algorithm, as well as the (information) complexity, crucially depend on the singular value decomposition of S_d . Hence, we have to calculate the eigenpairs $(\lambda_{d,i},\phi_{d,i})$ of the tensor product operator W_d obtained in (2.14). We follow [NW08, Section 5.2] and claim that these eigenpairs are given by (tensor) products of the non-trivial eigenpairs (λ_m,ϕ_m) , $m \in \mathcal{M}_1 = \{m \in \mathbb{N} \mid m < v(W_1) + 1\}$, of the univariate operator $W_1 = S_1^{\dagger}S_1$ (see (2.5)). This is the subject of the following assertion.

PROPOSITION 2.10. For $d \in \mathbb{N}$ the non-trivial eigenpairs of the operator $W_d = S_d^{\dagger} S_d$ are given by $\{(\widetilde{\lambda}_{d,\boldsymbol{m}}, \widetilde{\phi}_{d,\boldsymbol{m}}) \mid \boldsymbol{m} = (m_1, \dots, m_d) \in \mathcal{M}_d = (\mathcal{M}_1)^d\}$, where

$$\widetilde{\lambda}_{d,m} = \prod_{k=1}^{d} \lambda_{m_k} \quad and \quad \widetilde{\phi}_{d,m} = \bigotimes_{k=1}^{d} \phi_{m_k}.$$
 (2.15)

Proof. Obviously, all the $\widetilde{\phi}_{d,\boldsymbol{m}}$'s are mutually orthonormal in \mathcal{F}_d , i.e.

$$\langle \widetilde{\phi}_{d, \boldsymbol{i}}, \widetilde{\phi}_{d, \boldsymbol{j}} \rangle_{\mathcal{F}_d} = \prod_{k=1}^d \langle \phi_{i_k}, \phi_{j_k} \rangle_{\mathcal{F}_1} = \prod_{k=1}^d \delta_{i_k, j_k} = \delta_{\boldsymbol{i}, \boldsymbol{j}}, \quad \boldsymbol{i}, \boldsymbol{j} \in \mathcal{M}_d.$$

Furthermore,

$$W_{d}\widetilde{\phi}_{d,\boldsymbol{m}} = \left(\bigotimes_{k=1}^{d} W_{1}\right) \left(\bigotimes_{k=1}^{d} \phi_{m_{k}}\right) = \bigotimes_{k=1}^{d} \left(W_{1}\phi_{m_{k}}\right)$$
$$= \bigotimes_{k=1}^{d} \left(\lambda_{m_{k}}\phi_{m_{k}}\right) = \prod_{k=1}^{d} \lambda_{m_{k}} \cdot \bigotimes_{k=1}^{d} \phi_{m_{k}} = \widetilde{\lambda}_{d,\boldsymbol{m}}\widetilde{\phi}_{d,\boldsymbol{m}}$$

shows that $\widetilde{\phi}_{d,m}$, $m \in \mathcal{M}_d$, is indeed an eigenelement with respect to the strictly positive eigenvalue $\widetilde{\lambda}_{d,m}$ of W_d .

Assume for a moment there exists an eigenpair (μ, η) of W_d with $\mu \neq 0$ which cannot be represented by (2.15). Then, by the assertions in the previous section, η is orthogonal to any other eigenelement $\widetilde{\phi}_{d,\boldsymbol{m}}$, $\boldsymbol{m} \in \mathcal{M}_d$. Remember that $\Phi_1 = \{\phi_m \in \mathcal{F}_1 \mid m \in \mathcal{M}_1\}$ can be extended to an orthonormal basis $E_1 = \{e_m \mid m \in \mathcal{I}_1\}$ of \mathcal{F}_1 (see Remark 2.6) which can be used to construct an ONB $E_d = \{e_{d,\boldsymbol{j}} \mid \boldsymbol{j} \in \mathcal{I}_d = (\mathcal{I}_1)^d\}$ of \mathcal{F}_d given by (2.13). Therefore η can be represented as

$$\eta = \sum_{\boldsymbol{j} \in \mathcal{I}_d} \langle \eta, e_{d, \boldsymbol{j}} \rangle_{\mathcal{F}_d} e_{d, \boldsymbol{j}} = \sum_{\boldsymbol{j} \in \mathcal{M}_d} \langle \eta, \widetilde{\phi}_{d, \boldsymbol{j}} \rangle_{\mathcal{F}_d} \widetilde{\phi}_{d, \boldsymbol{j}} + \sum_{\boldsymbol{j} \in \mathcal{I}_d \setminus \mathcal{M}_d} \langle \eta, e_{d, \boldsymbol{j}} \rangle_{\mathcal{F}_d} e_{d, \boldsymbol{j}}
= \sum_{\boldsymbol{j} \in \mathcal{I}_d \setminus \mathcal{M}_d} \langle \eta, e_{d, \boldsymbol{j}} \rangle_{\mathcal{F}_d} e_{d, \boldsymbol{j}},$$

where each sum consists of at most countably many non-vanishing summands and converges unconditionally. Now the boundedness of S_d implies

$$S_d \eta = \sum_{\boldsymbol{j} \in \mathcal{I}_d \setminus \mathcal{M}_d} \langle \eta, e_{d, \boldsymbol{j}} \rangle_{\mathcal{F}_d} S_d e_{d, \boldsymbol{j}} = 0,$$

since each tensor product $S_d e_{d,j} = \bigotimes_{k=1}^d (S_1 e_{j_k})$, $j \in \mathcal{I}_d \setminus \mathcal{M}_d$, includes at least one factor $S_1 e_{j_k}$ with $j_k \notin \mathcal{M}_1$. These factors vanish because the set $\{S_1 e_m = S_1 \phi_m \mid m \in \mathcal{M}_1\}$ is an ONB of the image of S_1 in \mathcal{G}_1 . Hence, $W_d \eta = S_d^{\dagger}(S_d \eta) = 0$, which contradicts our assumption. In other words, (2.15) completely describes the eigenpairs of W_d , as claimed.

Again this proof justifies the restriction to separable spaces \mathcal{F}_1 (and hence also \mathcal{F}_d) in what follows (see Remark 2.6). Thus we can assume that the set of univariate eigenelements Φ_1 already is an ONB in \mathcal{F}_1 , i.e. that $\Phi_1 = E_1$, and consequently $\Phi_d = \{\widetilde{\phi}_{d,m} \mid m \in \mathcal{M}_d\}$ is an ONB in \mathcal{F}_d .

To unify our notation we rearrange the eigenpairs obtained according to the non-increasing ordering of the eigenvalues. To this end, note that $\#\mathcal{M}_d = (\#\mathcal{M}_1)^d$, i.e. we have $v(W_d) = v(W_1)^d$ strictly positive eigenvalues in dimension d. Therefore we define a sequence of bijections $\psi = \psi_d : \{i \in \mathbb{N} \mid i < v(W_1)^d + 1\} \to \mathcal{M}_d$ such that

$$\lambda_{d,i} = \widetilde{\lambda}_{d,\psi(i)} \ge \widetilde{\lambda}_{d,\psi(i+1)}$$
 for all $1 \le i < v(W_1)^d + 1$.

Consequently, the corresponding eigenelements are denoted by $\phi_{d,i} = \widetilde{\phi}_{d,\psi(i)}$. Similar to the definitions in Section 2.3.1 we extend the (possibly finite) sequence of eigenvalues by $\lambda_{d,i} = 0$ for $i > v(W_1)^d$. Observe that the largest eigenvalue in dimension d is given by

$$\lambda_{d,1} = \lambda_1^d$$

and thus the initial error is $\varepsilon_d^{\mathrm{init}} = \lambda_1^{d/2}$.

In virtue of Proposition 2.10, the optimal algorithm $A_{n,d}^*$ for linear tensor product problems $S = (S_d)_{d \in \mathbb{N}}$ can be deduced using Corollary 2.7. For $d \in \mathbb{N}$ and $n \in \mathbb{N}_0$ it reads

$$A_{n,d}^* \colon \mathcal{F}_d \to \mathcal{G}_d, \quad f \mapsto A_{n,d}^*(f) = \sum_{i=1}^{\min\{n, v(W_d)\}} \langle f, \phi_{d,i} \rangle_{\mathcal{F}_d} S_d \phi_{d,i}, \tag{2.16}$$

and its worst case error can be expressed in terms of the sequence $\lambda = (\lambda_m)_{m \in \mathbb{N}}$. More precisely, we have $e^{\text{wor}}(n,d;S_d) = \Delta^{\text{wor}}(A_{n,d}^*;S_d) = \sqrt{\lambda_{d,n+1}}$. We are ready to characterize tractability of such problems in the next subsection.

2.4.3. Complexity. We begin by analyzing the information complexity with respect to the absolute error criterion. Let $S_1 \colon \mathcal{F}_1 \to \mathcal{G}_1$ denote a compact linear operator between arbitrary Hilbert spaces \mathcal{F}_1 and \mathcal{G}_1 and let $S = (S_d)_{d \in \mathbb{N}}$ be the sequence of d-fold tensor product problems defined in Section 2.4.1. As before, the non-increasing sequence of nonnegative eigenvalues of the univariate operator $W_1 = S_1^{\dagger} S_1$ is denoted by $\lambda = (\lambda_m)_{m \in \mathbb{N}}$. At this point we stress that it is reasonable to assume that $\lambda_2 > 0$. Otherwise for every $d \in \mathbb{N}$ there would be only at most one non-vanishing d-dimensional eigenvalue of $W_d = S_d^{\dagger} S_d$. Hence the problem S_d would be trivial since then $n_{\text{abs}}^{\text{wor}}(\varepsilon, d) \leq 1$ for all $\varepsilon > 0$. Note that $\lambda_2 > 0$ also implies $\lambda_1 > 0$ so that S_1 and S_d are not the zero operator.

We now present an assertion which is mainly based on Theorem 5.5 in Novak and Woźniakowski [NW08]. The missing sufficient condition for weak tractability was later given by Papageorgiou and Petras [PP09]. Although the results of these authors only refer to linear tensor product problems defined between Hilbert function spaces they remain valid in our more general setting.

THEOREM 2.11. Consider the problem $S = (S_d)_{d \in \mathbb{N}}$ as described before. We study the absolute error criterion in the worst case setting.

- Let $\lambda_1 > 1$. Then S suffers from the curse of dimensionality.
- Let $\lambda_1 = 1$. Then
 - \circ S is polynomially intractable. In particular, if $\lambda_2 = 1$ then S suffers from the curse of dimensionality.
 - o S is weakly tractable if and only if $\lambda_2 < 1$ and $\lambda_n \in o(\ln^{-2} n)$ as $n \to \infty$ (12).
- Let $\lambda_1 < 1$. Then
 - S never suffers from the curse of dimensionality.
 - o S is weakly tractable if and only if $\lambda_n \in o(\ln^{-2} n)$ as $n \to \infty$.
 - o S is polynomially tractable if and only if it is strongly polynomially tractable. Moreover, this holds if and only if there exists some $\tau \in (0, \infty)$ such that $\lambda \in \ell_{\tau}$ and the exponent of strong polynomial tractability is given by

$$p^* = \inf \Big\{ 2\tau \; \Big| \; \sum_{m=1}^{\infty} \lambda_m^{\tau} \le 1 \Big\}.$$

For the sake of completeness we mention that Theorem 5.5 in [NW08] includes some additional lower bounds on the information complexity in the case $\lambda_1 \geq 1$. For poly-

⁽¹²⁾ To avoid possible confusions, here and in what follows, $\ln^{\alpha} n$ means $(\ln n)^{\alpha}$ where $\alpha \in \mathbb{R}$.

nomial (in)tractability the main idea of the proof is to apply Theorem 2.8 and to use the product structure of the sequences $(\lambda_{d,i})_{i\in\mathbb{N}}$ involved which are essentially given by Proposition 2.10. We will not provide an explicit proof here. Instead the interested reader is referred to Example 3.9 where we deduce all assertions stated in Theorem 2.11 from a generalized result for *scaled* tensor product problems. To get these more general assertions we will exactly follow the sketch of proof just mentioned.

Many authors in IBC use phrases like "(unweighted) tensor product problems are intractable". In this regard they refer to the following theorem for the normalized error criterion, which is essentially based on Theorem 5.6 of [NW08], as well as on [PP09]. From our point of view it is just a simple consequence of the assertions for absolute errors.

THEOREM 2.12. Consider the problem $S = (S_d)_{d \in \mathbb{N}}$ as described above. We study the normalized error criterion in the worst case setting.

- Let $\lambda_1 = \lambda_2$. Then S suffers from the curse of dimensionality.
- Let $\lambda_1 > \lambda_2$. Then
 - \circ S is weakly tractable if and only if $\lambda_n \in o(\ln^{-2} n)$ as $n \to \infty$.
 - \circ S is polynomially intractable.

Since the subsequent proof technique is typical in this field of research, we include the proof of Theorem 2.12 in full detail.

Proof. Assume we have already proven Theorem 2.11. Given the problem $S = (S_d)_{d \in \mathbb{N}}$, constructed out of $S_1 \colon \mathcal{F}_1 \to \mathcal{G}_1$, as well as the associated sequence $(\lambda_m)_{m \in \mathbb{N}}$, we define a new operator $T_1 \colon \mathcal{F}_1 \to \mathcal{G}_1$ by $f \mapsto T_1 f = (1/\sqrt{\lambda_1})S_1 f$. Clearly, T_1 is a linear and compact mapping between Hilbert spaces, and

$$\langle T_1 f, g \rangle_{\mathcal{G}_1} = \frac{1}{\sqrt{\lambda_1}} \langle S_1 f, g \rangle_{\mathcal{G}_1} = \frac{1}{\sqrt{\lambda_1}} \langle f, S_1^{\dagger} g \rangle_{\mathcal{F}_1} = \left\langle f, \left(\frac{1}{\sqrt{\lambda_1}} S_1^{\dagger} \right) g \right\rangle_{\mathcal{F}_1}$$

for $f \in \mathcal{F}_1$ and $g \in \mathcal{G}_1$. Hence, $T_1^{\dagger} = (1/\sqrt{\lambda_1})S_1^{\dagger}$ and the (extended) eigenvalue sequence of $V_1 = W_1(T) = T_1^{\dagger}T_1 = (1/\lambda_1)S_1^{\dagger}S_1 = (1/\lambda_1)W_1(S)$ is given by $\mu = (\mu_m)_{m \in \mathbb{N}}$, where $\mu_m = \lambda_m/\lambda_1$ for $m \in \mathbb{N}$. For details, see also the arguments used in Section 3.1. Anyway, using the mapping T_1 , we can construct the tensor product problem $T = (T_d)_{d \in \mathbb{N}}$ by the usual procedure. Now (2.15) in Proposition 2.10 shows that the corresponding eigenvalues of $V_d = W_d(T) = T_d^{\dagger}T_d$ are given by

$$\widetilde{\mu}_{d,\boldsymbol{m}} = \prod_{k=1}^{d} \mu_{m_k} = \frac{1}{\lambda_1^d} \prod_{k=1}^{d} \lambda_{m_k} = \frac{1}{\lambda_{d,1}} \widetilde{\lambda}_{d,\boldsymbol{m}}, \quad \boldsymbol{m} = (m_1, \dots, m_d) \in \mathcal{M}_d,$$

such that $\mu_{d,i} = (1/\lambda_{d,1})\lambda_{d,i}$ for $i \in \mathbb{N}$. This implies that the information complexity of S with respect to the normalized error criterion coincides with the absolute information complexity of the (scaled) problem T, i.e.

$$n_{\text{norm}}(\varepsilon', d; S_d) = \#\{n \in \mathbb{N}_0 \mid \lambda_{d,n}/\lambda_{d,1} > (\varepsilon')^2\} = \#\{n \in \mathbb{N}_0 \mid \mu_{d,n} > (\varepsilon')^2\}$$
$$= n_{\text{abs}}(\varepsilon', d; T_d)$$

for all $\varepsilon' \in (0,1]$ and each $d \in \mathbb{N}$. Since $\mu_1 = 1 \ge \mu_2 > 0$ we are allowed to apply Theorem 2.11 for T. Finally the observations that $\mu_2 = 1$ if and only if $\lambda_1 = \lambda_2$, as well

as that $\mu_n \in o(\ln^{-2} n)$ (as $n \to \infty$) if and only if $\lambda = (\lambda_n)_{n \in \mathbb{N}}$ belongs to this class, complete the proof.

2.5. Reproducing kernel Hilbert spaces. When we deal with problems defined on Hilbert function spaces H, a special kind of Hilbert spaces is of particular interest. The reason is that in practice often only function evaluations rather than information obtained by arbitrary linear functionals are permitted. In order to compare the power of these classes of information operations (Λ^{std} vs. Λ^{all}) from a theoretical point of view, it seems useful to investigate conditions which ensure that point evaluation functionals

$$L_y \colon H \to \mathbb{R}, \quad f \mapsto L_y(f) = f(y),$$

for all y in the domain of definition Ω of $f \in H$, belong to the class $\Lambda^{\rm all}$. Clearly L_y is always linear so it is enough to ask whether it is also continuous (or bounded, respectively) in f. It turns out that, as long as we restrict ourselves to Hilbert spaces, this property can be characterized by the existence of a so-called reproducing kernel K. If so, then the space H is referred to as a reproducing kernel Hilbert space (RKHS for short) and we write $H = \mathcal{H}(K)$. In the present section we collect some basic properties of this concept which we will need later on in Section 4.1.2. The presentation given here is based on the famous paper of Aronszajn [Aro50], as well as the textbook of Wahba [Wah90] (13). Standard examples of RKHSs, such as Korobov spaces and Sobolev spaces of dominating mixed smoothness, can be found in [NW08, Appendix A].

- **2.5.1. Definition and properties.** A (real) Hilbert space H of functions $f: \Omega \to \mathbb{R}$, equipped with inner product $\langle \cdot, \cdot \rangle_H$, is said to be a reproducing kernel Hilbert space if there exists a function $K: \Omega \times \Omega \to \mathbb{R}$ such that
- for all fixed $y \in \Omega$ the function $K_y = K(\cdot, y)$ belongs to H, and
- for every $f \in H$ and all $y \in \Omega$,

$$L_y(f) = f(y) = \langle f, K_y \rangle_H = \langle f, K(\cdot, y) \rangle_H. \tag{2.17}$$

Formula (2.17) is known as the reproducing property. Together with the first point it obviously implies the boundedness of point evaluations on $H = \mathcal{H}(K)$. The converse, i.e. the existence (and uniqueness) of the reproducing kernel K, is a simple consequence of the Riesz representation theorem (see [Tri92, p. 90] or [Yos80, III.6]). Unfortunately the proof of this theorem is non-constructive and therefore it does not provide an explicit method to find the representer $K_y = K(\cdot, y)$ of L_y . In fact, given a specific RKHS $\mathcal{H}(K)$ it seems to be a challenging problem to deduce a closed form of its reproducing kernel K. However, as long as we restrict ourselves to separable RKHSs, it is easy to prove that K is given by

$$K(x,y) = \sum_{m \in \mathcal{I}} e_m(x) e_m(y), \quad x, y \in \Omega,$$
(2.18)

where $\{e_m : \Omega \to \mathbb{R} \mid m \in \mathcal{I}\}$ denotes an arbitrary orthonormal basis of $\mathcal{H}(K)$. Furthermore we know that every reproducing kernel K is positive definite. That is, for all $n \in \mathbb{N}$

 $[\]binom{13}{}$ For the ease of notation (and in contrast to the above mentioned references) we restrict ourselves to spaces over \mathbb{R} . Once more the theory can be transferred almost literally to \mathbb{C} .

and any sequence $\mathbf{x} = (x_m)_{m=1}^n \in \Omega^n$ the quadratic form

$$Q_{K;x}(\xi_1,\ldots,\xi_n) = \sum_{i,j=1}^n K(x_i,x_j)\,\xi_i\,\xi_j, \quad \boldsymbol{\xi} = (\xi_m)_{m=1}^n \in \mathbb{R}^n, \quad (2.19)$$

is a non-negative function of ξ . In particular,

$$K(x,x) \ge 0$$
 and $K(x,y) = K(y,x)$, for all $x,y \in \Omega$.

Conversely, Moore showed that every positive definite function K in the above sense uniquely determines a RKHS H admitting K as its reproducing kernel; see [Aro50]. Again it turned out to be a hard problem to determine a suitable representation of H (and its inner product) for a given function K.

Besides further fascinating properties, we want to focus on products of kernel functions. To this end, for $d \in \mathbb{N}$ let $K^{(k)}$, $k = 1, \ldots, d$, denote a finite number of reproducing kernels defined on the sets $\Omega^{(k)} \times \Omega^{(k)}$, respectively. Then we may consider the tensor product

$$K_d = \bigotimes_{k=1}^d K^{(k)} : \left(\left. \bigotimes_{k=1}^d \Omega^{(k)} \right) \times \left(\left. \bigotimes_{k=1}^d \Omega^{(k)} \right) \to \mathbb{R}, \quad K_d(\boldsymbol{x}, \boldsymbol{y}) = \prod_{k=1}^d K^{(k)}(x_k, y_k), \right.$$

where we set $\mathbf{x} = (x_1, \dots, x_d)$ and $\mathbf{y} = (y_1, \dots, y_d)$ with $x_k, y_k \in \Omega^{(k)}$. On the other hand, each kernel $K^{(k)}$ induces a uniquely defined RKHS $\mathcal{H}(K^{(k)})$ which in turn implies the existence of one (and only one) tensor product space $H_d = \bigotimes_{k=1}^d \mathcal{H}(K^{(k)})$ using the arguments presented in Section 2.4. Now it can be checked that H_d itself is a RKHS and its kernel is given by K_d , i.e.

$$H_d = \mathcal{H}(K_d) = \bigotimes_{k=1}^d \mathcal{H}(K^{(k)}).$$
 (2.20)

The proof of this assertion can be obtained inductively by adding one factor in every step. Then it remains to show that the resulting quadratic forms (2.19) are non-negative again, which can be done using a classical result due to Schur.

Note that the whole theory works for arbitrary point sets Ω ; this turned out to be useful in the context of so-called support vector machines, which are instances of the more general class of kernel methods. However, in IBC special choices such as $\Omega = \Omega_1 = [0,1]$ (or $\Omega = \mathbb{R}$) are of particular interest. For multivariate problems the standard choice is $\Omega = \Omega_d = \Omega_1^d$, which perfectly fits the tensor product construction explained before. In this respect the univariate kernels $K^{(k)}$, $k = 1, \ldots, d$, are often taken as weighted instances $K_1^{\gamma_{d,k}}$ of some underlying kernel K_1 . A prominent example is given by $K_1^{\gamma_{d,k}}(x,y) = 1 + \gamma_{d,k} \min\{x,y\}$, which leads to an anchored Sobolev space $\widetilde{\mathcal{H}}_d^{\gamma}$ related to the Wiener sheet measure (see, e.g., [KWW08, Section 8] or [Wei12b]). Another example of this type will be discussed in detail within Section 4.1.2.

Finally we mention that the concept of RKHSs was generalized recently to the class of so-called *reproducing kernel Banach spaces* (*RKBSs*). For a brief introduction to this topic we refer to Zhang and Zhang [ZZ13].

2.5.2. Examples: Integration and approximation problems. Let us conclude the presentation with some examples which show that the knowledge about the existence of

a reproducing kernel can be exploited to obtain complexity assertions for the classical problems of integration and approximation.

EXAMPLE 2.13 (Worst case error of QMC rules). For $d \in \mathbb{N}$ suppose $\mathcal{H}(K_d)$ is a RKHS of real-valued functions f defined on some Borel measurable subset Ω_d of \mathbb{R}^d . Consider the solution operator of the integration problem

$$\operatorname{Int}_d^{\varrho_d} \colon \mathcal{B}(\mathcal{H}(K_d)) o \mathbb{R}, \quad f \mapsto \operatorname{Int}_d^{\varrho_d} f = \int_{\Omega_d} f(\boldsymbol{x}) \, \varrho_d(\boldsymbol{x}) \, \mathrm{d}\lambda^d(\boldsymbol{x}),$$

where ϱ_d denotes a probability density function on Ω_d . Let us additionally assume that the function

$$h_d = \int_{\Omega_d} K_d(\cdot, \boldsymbol{x}) \varrho_d(\boldsymbol{x}) \, \mathrm{d}\lambda^d(\boldsymbol{x})$$

is well-defined and belongs to $\mathcal{H}(K_d)$. Then it is easy to see that h_d is the representer of the linear functional $\mathrm{Int}_d^{\varrho_d}$, i.e. that $\mathrm{Int}_d^{\varrho_d} f = \langle f, h_d \rangle_{\mathcal{H}(K_d)} < \infty$ for all $f \in \mathcal{H}(K_d)$. Since allowing arbitrary linear functionals to approximate the value of the integral would make the problem trivial we consider cubature rules of the form

$$A_{n,d}f = \sum_{i=1}^{n} a_i f(\boldsymbol{x}^{(i)}), \quad n \in \mathbb{N}_0,$$

defined by a priori chosen sample points $\mathbf{x}^{(i)} \in \Omega_d$ and some weights $a_i \in \mathbb{R}$, i = 1, ..., n. Due to (2.17) the linear operator $A_{n,d}$ possesses a representer in the space $\mathcal{H}(K_d)$ as well. Consequently, its worst case error can be computed exactly in terms of the reproducing kernel and the parameters $(a_i)_{i=1}^n$ and $(\mathbf{x}^{(i)})_{i=1}^n$:

$$\Delta^{\text{wor}}(A_{n,d}; \operatorname{Int}_{d}^{\varrho_{d}} : \mathcal{B}(\mathcal{H}(K_{d})) \to \mathbb{R})^{2}$$

$$= \sup_{f \in \mathcal{B}(\mathcal{H}(K_{d}))} \left| \left(\operatorname{Int}_{d}^{\varrho_{d}} - A_{n,d} \right) (f) \right|^{2} = \left\| h_{d} - \sum_{i=1}^{n} a_{i} K_{d}(\cdot, \boldsymbol{x}^{(i)}) \mid \mathcal{H}(K_{d}) \right\|^{2}$$

$$= \int_{\Omega_{d}^{2}} K_{d}(\boldsymbol{x}, \boldsymbol{y}) \varrho_{d}(\boldsymbol{x}) \varrho_{d}(\boldsymbol{y}) d\lambda^{2d}(\boldsymbol{x}, \boldsymbol{y}) - 2 \sum_{i=1}^{n} a_{i} \int_{\Omega_{d}} K_{d}(\boldsymbol{x}, \boldsymbol{x}^{(i)}) \varrho_{d}(\boldsymbol{x}) d\lambda^{d}(\boldsymbol{x})$$

$$+ \sum_{i, i=1}^{n} a_{i} a_{j} K_{d}(\boldsymbol{x}^{(i)}, \boldsymbol{x}^{(j)}).$$

Choosing special weights a_i (such as $a_i \equiv 1/n$), as well as specific sample points $\boldsymbol{x}^{(i)}$ (e.g. points from a so-called *integration lattice*), we end up with well-studied classes of cubature rules which are known as *quasi-Monte Carlo* (QMC) methods and *lattice rules*, respectively. The common feature of those integration schemes is that their complexity analysis is mainly based on the presented worst case error formula and thus on the properties of the reproducing kernel K_d . Moreover, the above expression for $\Delta^{\text{wor}}(A_{n,d})$ plays an important role in *discrepancy theory*.

Various kinds of integration problems are studied in Novak and Woźniakowski [NW10]. For the recent state of the art in discrepancy theory and QMC methods we refer the reader to the monograph of Dick and Pillichshammer [DP10], as well as to the survey article of

Dick, Kuo and Sloan [DKS13] and the references therein. An introduction to lattice rules can also be found in the textbook of Sloan and Joe [SJ94]. \Box

Our second example shows the relation of reproducing kernels and the singular values for certain approximation operators.

EXAMPLE 2.14 (Weighted L₂-approximation). For $d \in \mathbb{N}$ let $\mathcal{H}(K_d)$ be a separable and infinite-dimensional RKHS which is compactly embedded into $L_2^{\varrho_d}(\Omega_d)$. Here ϱ_d again denotes some probability density on $\Omega_d \subseteq \mathbb{R}^d$. Then we may study the approximation problem

$$\operatorname{App}_{d}^{\varrho_{d}} : \mathcal{B}(\mathcal{H}(K_{d})) \to \operatorname{L}_{2}^{\varrho_{d}}(\Omega_{d}), \quad f \mapsto \operatorname{App}_{d}^{\varrho_{d}} f = f, \tag{2.21}$$

in the worst case setting. Since both the source and target spaces are Hilbert spaces we can use the theory developed in Section 2.3 to obtain complexity results with respect to the class $\Lambda^{\rm all}$. Therefore we need to analyze the eigenvalues of the compact operator $W_d^{\varrho_d} = ({\rm App}_d^{\varrho_d})^{\dagger} {\rm App}_d^{\varrho_d}$. Using the reproducing property (2.17) and the symmetry of K_d we conclude that

$$\begin{split} \left(W_d^{\varrho_d} f\right)(\boldsymbol{x}) &= \langle (\operatorname{App}_d^{\varrho_d})^{\dagger} \operatorname{App}_d^{\varrho_d} f, K_d(\cdot, \boldsymbol{x}) \rangle_{\mathcal{H}(K_d)} = \langle \operatorname{App}_d^{\varrho_d} f, \operatorname{App}_d^{\varrho_d} K_d(\cdot, \boldsymbol{x}) \rangle_{\operatorname{L}_2^{\varrho_d}(\Omega_d)} \\ &= \int_{\Omega_d} f(\boldsymbol{y}) \, K_d(\boldsymbol{x}, \boldsymbol{y}) \, \varrho_d(\boldsymbol{y}) \, \mathrm{d}\lambda^d(\boldsymbol{y}) \end{split}$$

for all $f \in \mathcal{H}(K_d)$ and any $\boldsymbol{x} \in \Omega_d$. Hence, $W_d^{\varrho_d}$ takes the form of a weighted integral operator against the kernel $K_d(\cdot, \boldsymbol{y})$ and its non-trivial eigenpairs $\{(\lambda_{d,\varrho_d,i}, \phi_{d,\varrho_d,i}) \mid i \in \mathcal{M}_d\}$ can be found by solving integral equations. Formula (2.18) yields

$$K_d(\boldsymbol{x}, \boldsymbol{x}) = \sum_{i \in \mathcal{M}_d} \phi_{d, \varrho_d, i}(\boldsymbol{x})^2 < \infty$$

for every $\boldsymbol{x} \in \Omega_d$ because we know that $\{\phi_{d,\varrho_d,i} \mid i \in \mathcal{M}_d\}$ forms an ONB in $\mathcal{H}(K_d)$. Since $\phi_{d,\varrho_d,i} \in L_2^{\varrho_d}(\Omega_d)$ and $\|\phi_{d,\varrho_d,i} \mid L_2^{\varrho_d}(\Omega_d)\|^2 = \lambda_{d,\varrho_d,i}$ for all $i \in \mathcal{M}_d$, it easily follows that

trace
$$W_d^{\varrho_d} = \sum_{i \in \mathcal{M}_d} \lambda_{d,\varrho_d,i} = \sum_{i \in \mathcal{M}_d} \|\phi_{d,\varrho_d,i} \mid \mathcal{L}_2^{\varrho_d}(\Omega_d)\|^2$$

$$= \int_{\Omega_d} K_d(\boldsymbol{x}, \boldsymbol{x}) \,\varrho_d(\boldsymbol{x}) \,\mathrm{d}\lambda^d(\boldsymbol{x}). \tag{2.22}$$

Note that this trace may be finite or infinite depending on the values of K_d on the diagonal $\{(x,x) \mid x \in \Omega_d\}$. It turns out that an infinite trace implies that there is, in general, no (non-trivial) relation of the power of Λ^{all} and Λ^{std} for the given approximation problem. In contrast, it is known that for finite traces there exist close relations of these classes of information operations. In particular, it is possible to derive bounds on the rate of convergence for Λ^{std} from corresponding bounds for Λ^{all} . For details we refer to [NW12, Chapter 26].

Finally we note that the finite trace property of $W_d^{\varrho_d}$ immediately implies that $\lambda_{d,\varrho_d,i} \in \mathcal{O}(i^{-1})$ as $i \to \infty$. Hence, if we deal with linear information, we can deduce that $e^{\mathrm{wor}}(n,d;\mathrm{App}_d^{\varrho_d}) \in \mathcal{O}(n^{-1/2}), n \to \infty$, directly from an integrability property of the kernel K_d . \square

In the last example we present a useful relation of reproducing kernels and average case approximation problems.

EXAMPLE 2.15 (Average case approximation). For $d \in \mathbb{N}$ assume ϱ_d is some probability density function on $\Omega_d = [0,1]^d$ and let $K_d \colon \Omega_d \times \Omega_d \to \mathbb{R}$ denote a reproducing kernel such that the mapping $\boldsymbol{x} \mapsto K_d(\boldsymbol{x}, \boldsymbol{x})$ belongs to $L_1^{\varrho_d}(\Omega_d)$. That is, suppose (2.22) is finite. Furthermore, let \mathcal{F}_d denote a separable Banach space of real-valued functions on Ω_d which is continuously embedded into $L_2^{\varrho_d}(\Omega_d)$ and for which function evaluations are continuous. We equip \mathcal{F}_d with a zero-mean Gaussian measure μ_d such that its correlation operator $C_{\mu_d} \colon \mathcal{F}_d^* \to \mathcal{F}_d$ applied to point evaluation functionals $L_{\boldsymbol{x}}$ can be expressed in terms of K_d :

$$K_d(\boldsymbol{x}, \boldsymbol{y}) = L_{\boldsymbol{x}}(C_{\mu_d}L_{\boldsymbol{y}}) = \int_{\mathcal{F}_d} f(\boldsymbol{x}) f(\boldsymbol{y}) d\mu_d(f)$$
 for all $\boldsymbol{x}, \boldsymbol{y} \in \Omega_d$.

We stress that this is always possible for a suitable choice of \mathcal{F}_d and that our assumptions imply a continuous embedding of the RKHS $\mathcal{H}(K_d)$ (induced by K_d) into \mathcal{F}_d . Consequently, also $\operatorname{App}_d^{\varrho_d} \colon \mathcal{H}(K_d) \to \operatorname{L}_2^{\varrho_d}(\Omega_d)$ is bounded, i.e. continuous (see (2.21)). For details and concrete examples the reader is referred to [NW08, Appendix B], [NW10, Section 13.2], and [NW12, Section 24.1].

As in the previous example we want to look for good approximations $A_{n,d}f$ to input functions f in the norm of $L_2^{\varrho_d}(\Omega_d)$. This time we measure the average performance of the algorithm $A_{n,d}$ with respect to μ_d , i.e. we try to minimize

$$\Delta^{\operatorname{avg}}\left(A_{n,d};\operatorname{id}_{d}^{\varrho_{d}}\colon\mathcal{F}_{d}\to\operatorname{L}_{2}^{\varrho_{d}}(\Omega_{d})\right)=\left(\int_{\mathcal{F}_{d}}\|f-A_{n,d}f\mid\operatorname{L}_{2}^{\varrho_{d}}(\Omega_{d})\|^{2}\operatorname{d}\mu_{d}(f)\right)^{1/2}.$$

Observe that $\nu_d = \mu_d \circ (\mathrm{id}_d^{\varrho_d})^{-1}$ defines a Gaussian measure on the subset $\mathrm{id}_d^{\varrho_d}(\mathcal{F}_d)$ of $L_2^{\varrho_d}(\Omega_d)$. Now it can be checked that the corresponding covariance operator

$$C_{\nu_d}^{\varrho_d} \colon \operatorname{id}_d^{\varrho_d}(\mathcal{F}_d) \to L_2^{\varrho_d}(\Omega_d)$$

of the measure ν_d is given by

$$f \mapsto (C_{\nu_d}^{\varrho_d} f)(\cdot) = \int_{\Omega_d} f(\boldsymbol{y}) K_d(\cdot, \boldsymbol{y}) \varrho_d(\boldsymbol{y}) d\lambda^d(\boldsymbol{y}).$$

This operator is self-adjoint, compact and has a finite trace due to the integrability assumption on K_d . Consequently, there exists a countable set of non-trivial eigenpairs $(\lambda_{d,\varrho_d,i},\eta_{d,\varrho_d,i})$ where the eigenfunctions $\eta_{d,\varrho_d,i}$ are mutually orthogonal (and normalized) with respect to the $L_2^{\varrho_d}(\Omega_d)$ -norm (see also Hickernell and Woźniakowski [HW00]).

Once more it turns out that the optimal algorithm $A_{n,d}^*$ in this setting is given by the orthogonal projection of the input function onto the subspace spanned by the eigenfunctions $\eta_{d,\varrho_d,i}$ which correspond to the n largest eigenvalues $\lambda_{d,\varrho_d,i}$. In contrast to the worst case setting the nth minimal average case error is

$$e^{\operatorname{avg}}(n,d;\operatorname{id}_d^{\varrho_d}:\mathcal{F}_d\to \operatorname{L}_2^{\varrho_d}(\Omega_d))=\Big(\sum_{i=n+1}^{\infty}\lambda_{d,\varrho_d,i}\Big)^{1/2},\quad n\in\mathbb{N}_0,\,d\in\mathbb{N},$$

if we assume a non-increasing ordering of the sequence $(\lambda_{d,\varrho_d,i})_{i=1}^{\infty}$ (14). Based on the

 $^(^{14})$ For the ease of notation we assumed here that all the eigenvalues are strictly positive.

above error formula it is possible to obtain characterizations of several types of tractability similar to the assertions given in Section 2.4.3 (see, e.g., [NW08, Chapter 6]).

We complete the discussion with the observation that the sets of (non-trivial) eigenpairs $(\lambda_{d,\varrho_d,i},\eta_{d,\varrho_d,i})$ of the operators $C^{\varrho_d}_{\nu_d}$ as defined above and $W^{\varrho_d}_d$ from Example 2.14 coincide, since $C^{\varrho_d}_{\nu_d}$ only takes values in $\mathcal{H}(K_d)$. To be precise, we note that $K_d(\boldsymbol{x},\boldsymbol{y}) = K_d(\boldsymbol{y},\boldsymbol{x})$ equals $(\mathrm{App}_d^{\varrho_d}K_d(\cdot,\boldsymbol{x}))(\boldsymbol{y})$ for each fixed \boldsymbol{x} and λ^d -almost every $\boldsymbol{y} \in \Omega_d$. Hence the chain of equations

$$(C_{\nu_d}^{\varrho_d} f)(\boldsymbol{x}) = \int_{\Omega_d} f(\boldsymbol{y}) \left(\operatorname{App}_d^{\varrho_d} K_d(\cdot, \boldsymbol{x}) \right) (\boldsymbol{y}) \, \varrho_d(\boldsymbol{y}) \, \mathrm{d}\lambda^d(\boldsymbol{y}) = \langle f, \operatorname{App}_d^{\varrho_d} K_d(\cdot, \boldsymbol{x}) \rangle_{\operatorname{L}_2^{\varrho_d}(\Omega_d)}$$
$$= \langle (\operatorname{App}_d^{\varrho_d})^{\dagger} f, K_d(\cdot, \boldsymbol{x}) \rangle_{\mathcal{H}(K_d)} = ((\operatorname{App}_d^{\varrho_d})^{\dagger} f)(\boldsymbol{x})$$

holds true for every $f \in \mathrm{id}_d^{\varrho_d}(\mathcal{F}_d) \subset \mathrm{L}_2^{\varrho_d}(\Omega_d)$ and λ^d -almost all $\boldsymbol{x} \in \Omega_d$ (15). Clearly $(\mathrm{App}_d^{\varrho_d})^{\dagger}$ maps into $\mathcal{H}(K_d)$ per definition. Thus, for every eigenfunction $\eta \in \mathrm{id}_d^{\varrho_d}(\mathcal{F})$ of $C_{\nu_d}^{\varrho_d}$, i.e.

$$\lambda \eta = C_{\nu_d}^{\varrho_d} \eta = (\text{App}_d^{\varrho_d})^{\dagger} \eta, \quad \lambda^d$$
-a.e. on Ω_d , (2.23)

we can find a representer $\overline{\eta} \in \mathcal{H}(K_d)$ with $\overline{\eta} = \eta$ in the sense of $L_2^{\varrho_d}(\Omega_d)$, such that the equalities in (2.23) hold pointwise on the whole set Ω_d and therefore also in the norm of $\mathcal{H}(K_d)$. Now it is easy to check that $(\lambda, \overline{\eta})$ is indeed an eigenpair of $W_d^{\varrho_d} = (\mathrm{App}_d^{\varrho_d})^{\dagger} \mathrm{App}_d^{\varrho_d}$, normalized with respect to the $L_2^{\varrho_d}(\Omega_d)$ -norm. Conversely, every eigenpair $(\lambda, \overline{\eta})$ of the operator $W_d^{\varrho_d}$ obviously fulfills $\lambda \overline{\eta} = C_{\nu_d}^{\varrho_d} \overline{\eta}$ interpreted in $L_2^{\varrho_d}(\Omega_d)$.

In conclusion we see that the knowledge of these eigenpairs implies complexity assertions for both approximation problems in the respective (quite different) settings. \Box

⁽¹⁵⁾ Note that $C_{\nu_d}^{\varrho_d} f \in L_2^{\varrho_d}(\Omega_d)$, i.e. it is uniquely defined on Ω_d up to a set of measure zero.

3. Problems on Hilbert spaces with scaled norms

The present chapter deals with a generalization of tensor product problems $S = (S_d)_{d \in \mathbb{N}}$ between Hilbert spaces in the sense of Section 2.4. We introduce additional scaling factors s_d to the norm of the source spaces \mathcal{F}_d and analyze their influence on the squared singular values $\lambda_{d,s_d,i}$ of the new problem operators S_{d,s_d} . Using the techniques from Section 2.4.2 we conclude optimal algorithms for these modified problems at the end of Section 3.1. Afterwards, in Section 3.2, we investigate tractability properties of this class of problems with respect to the worst case setting. Finally we present some applications of our results in Section 3.3.

3.1. Definitions, eigenpairs and the optimal algorithm. Let H_1 and \mathcal{G}_1 be arbitrary Hilbert spaces with inner products $\langle \cdot, \cdot \rangle_{H_1}$ and $\langle \cdot, \cdot \rangle_{\mathcal{G}_1}$, respectively. Further assume $S_1 \in \mathcal{K}(H_1, \mathcal{G}_1)$ is a compact linear operator between these spaces. Following the constructions given in Section 2.4.1 for any $d \in \mathbb{N}$ there exist uniquely defined d-fold tensor product spaces of H_1 and \mathcal{G}_1 . Let us denote these spaces by $H_d = H_1 \otimes \cdots \otimes H_1$ and $\mathcal{G}_d = \mathcal{G}_1 \otimes \cdots \otimes \mathcal{G}_1$, respectively. Finally we define $S = (S_d)_{d \in \mathbb{N}}$ to be the sequence of multivariate tensor product operators constructed from S_1 .

In contrast to Section 2.4 we now adapt the source spaces of our multivariate problem by introducing an additional positive sequence of scaling factors $s = (s_d)_{d \in \mathbb{N}}$. That is, for every $d \in \mathbb{N}$ we define \mathcal{F}_d to be Hilbert space H_d equipped with the inner product

$$\langle \cdot, \cdot \rangle_{\mathcal{F}_d} = \frac{1}{s_d} \langle \cdot, \cdot \rangle_{H_d}, \quad \text{where} \quad s_d > 0.$$
 (3.1)

Obviously \mathcal{F}_d algebraically coincides with H_d whereas the norms (induced by the respective inner products) are equivalent. Accordingly, the operators S_d are still well-defined for any $d \in \mathbb{N}$ when we replace H_d by \mathcal{F}_d . On the other hand the approximability properties of S crucially depend on the norms used, since we need to consider the whole unit ball $\widetilde{\mathcal{F}}_d = \mathcal{B}(\mathcal{F}_d)$ when dealing with the worst case setting. So let us denote the modified problem by $S_{(s)} = (S_{d,s_d} \colon \mathcal{F}_d \to \mathcal{G}_d)_{d \in \mathbb{N}}$.

From Section 2.3 we know that for the nth optimal algorithms for $S_{(s)}$ we need to study the eigenpairs of $W_{d,s_d} = S_{d,s_d}^{\dagger} S_{d,s_d}$. Although S_{d,s_d} equals S_d (as a mapping) we cannot claim that $W_{d,s_d} = W_d$ since S_{d,s_d}^{\dagger} does not necessarily coincide with S_d^{\dagger} . Nevertheless, there exists a strong relation. The following proposition extends Proposition 2.10 to the case of scaled problems in the above mentioned sense. Keep in mind that the eigenpairs of the univariate (unscaled) operator $W_1 = S_1^{\dagger} S_1$ are given by $\{(\lambda_m, e_m) \mid m \in \mathcal{M}_1\}$, where $\mathcal{M}_1 = \{m \in \mathbb{N} \mid m < v(W_1) + 1\}$ and $0 < \lambda_{m+1} \le \lambda_m$ for all $m < v(W_1)$.

PROPOSITION 3.1. For $d \in \mathbb{N}$ the non-trivial eigenpairs of $W_{d,s_d} = S_{d,s_d}^{\dagger} S_{d,s_d}$ are given by $\{(\widetilde{\lambda}_{d,s_d,m}, \widetilde{\phi}_{d,s_d,m}) \mid m \in \mathcal{M}_d = (\mathcal{M}_1)^d\}$, where

$$\widetilde{\lambda}_{d,s_d,m} = s_d \, \widetilde{\lambda}_{d,m} = s_d \, \prod_{k=1}^d \lambda_{m_k} \quad and \quad \widetilde{\phi}_{d,s_d,m} = \sqrt{s_d} \, \widetilde{\phi}_{d,m} = \sqrt{s_d} \, \bigotimes_{k=1}^d \phi_{m_k}. \quad (3.2)$$

Proof. Since $S_{d,s_d}f = S_df$ for every $f \in \mathcal{F}_d$ (or H_d , respectively) we have

$$\langle S_{d,s_d}f,g\rangle_{\mathcal{G}_d} = \langle S_df,g\rangle_{\mathcal{G}_d} = \langle f,S_d^\dagger g\rangle_{H_d} = s_d\langle f,S_d^\dagger g\rangle_{\mathcal{F}_d} = \langle f,(s_dS_d^\dagger)g\rangle_{\mathcal{F}_d}$$

for all $f \in \mathcal{F}_d$ and $g \in \mathcal{G}_d$. Thus, (2.4) and the uniqueness of the adjoint operator (¹⁶) imply that $S_{d,s_d}^{\dagger} = s_d \cdot S_d^{\dagger}$ pointwise, and consequently W_{d,s_d} equals $s_d W_d$ as a mapping. Hence, from Proposition 2.10 and the linearity of W_d we conclude that (3.2) indeed are eigenpairs of W_{d,s_d} . Due to the factor $\sqrt{s_d}$ and the relation (3.1) the eigenelements $\widetilde{\phi}_{d,s_d,m}$ are properly normalized in \mathcal{F}_d .

It remains to show that there cannot exist eigenpairs other than (3.2). This can be seen using arguments similar to those in the second part of the proof of Proposition 2.10. To this end, note that by (3.1) the inner product in \mathcal{F}_d equals zero if and only if the elements under consideration are orthogonal in H_d .

Given Proposition 3.1, the rest of this section is straightforward. Namely, we can use the bijections $\psi = \psi_d$ from Section 2.4.2 to define the non-increasing sequences $(\lambda_{d,s_d,i})_{i\in\mathbb{N}}$ by

$$\lambda_{d,s_d,i} = \begin{cases} \widetilde{\lambda}_{d,s_d,\psi(i)} & \text{if } 1 \le i < v(W_1)^d + 1, \\ 0 & \text{otherwise,} \end{cases}$$

for every $d \in \mathbb{N}$. The corresponding reordered eigenelements are denoted by $\phi_{d,s_d,i}$ with $1 \leq i < v(W_1)^d + 1$. Moreover, we again use Corollary 2.7 to see that for any $d \in \mathbb{N}$ the nth optimal algorithm for S_{d,s_d} , $n \in \mathbb{N}_0$, is given by

$$A_{n,d,s_d}^* \colon \mathcal{F}_d \to \mathcal{G}_d, \quad f \mapsto A_{n,d,s_d}^*(f) = \sum_{i=1}^{\min\{n,v(W_1)^d\}} \langle f, \phi_{d,s_d,i} \rangle_{\mathcal{F}_d} S_{d,s_d} \phi_{d,s_d,i}.$$

It realizes the nth minimal worst case error in dimension d, which equals

$$e^{\text{wor}}(n, d; S_{d,s_d}) = \Delta^{\text{wor}}(A_{n,d,s_d}^*; S_{d,s_d}) = \sqrt{\lambda_{d,s_d,n+1}}.$$
 (3.3)

In particular, the case n = 0, i.e. the initial error

$$\varepsilon_d^{\text{init}} = \sqrt{\lambda_{d,s_d,1}} = \sqrt{s_d \lambda_1^d},$$

will play an important role in what follows.

3.2. Complexity. As in Section 2.4, we proceed with the analysis of the information complexity of scaled tensor product problems $S_{(s)} = (S_{d,s_d})_{d \in \mathbb{N}}$ in the worst case setting. We first take a look at necessary and sufficient conditions for (strong) polynomial tractability with respect to absolute errors. Afterwards, in Section 3.2.2, we complete

⁽¹⁶⁾ Note that, clearly, S_{d,s_d} is compact if and only if S_d is compact.

these assertions and investigate conditions for weak tractability and the curse of dimensionality. Finally we will see in Section 3.2.3 that the improvements obtained by scaling are completely cancelled when we turn to the normalized error criterion.

As usual $\lambda = (\lambda_m)_{m \in \mathbb{N}}$ denotes the (extended) sequence of squared singular values of the underlying operator $S_1 \colon H_1 \to \mathcal{G}_1$. To avoid triviality we assume that $\lambda_2 > 0$ throughout the rest of this section. The reason for this assumption is explicitly stated in Section 2.4.3.

3.2.1. Polynomial tractability. The next statement is originally based on Theorem 3.1 of Woźniakowski [Woź94b], which provided the underlying idea for [NW08, Theorem 5.5]. We extend the results stated there to the case of scaled problems.

THEOREM 3.2. Let $S_{(s)} = (S_{d,s_d})_{d \in \mathbb{N}}$ denote a scaled tensor product problem in the sense of Section 3.1. Assume $\lambda_2 > 0$ and consider the worst case setting with respect to the absolute error criterion. Then the following assertions are equivalent:

- (I) $S_{(s)}$ is strongly polynomially tractable.
- (II) $S_{(s)}$ is polynomially tractable.
- (III) There exists $\tau \in (0, \infty)$ such that $\lambda \in \ell_{\tau}$ and $\sup_{d \in \mathbb{N}} s_d \|\lambda\| \ell_{\tau} \|^d < \infty$.
- (IV) There exists $\varrho \in (0, \infty)$ such that $\lambda \in \ell_{\varrho}$ and $\limsup_{d \to \infty} s_d^{1/d} < 1/\lambda_1$.

If one (and hence all) of these conditions applies then the exponent of strong polynomial tractability is given by

$$p^* = \inf\{2\tau \mid \tau \text{ fulfills condition (III)}\}.$$

Proof. Step 1. Since (I) clearly implies (II) we start by proving (II) \Rightarrow (III). Therefore let $S_{(s)}$ be polynomially tractable with the constants C, p > 0 and $q \ge 0$. Then Theorem 2.8 shows, for all $\varrho > p/2$, that

$$0 < C_{\varrho} = \sup_{d \in \mathbb{N}} \frac{1}{d^{2q/p}} \Big(\sum_{i=\lceil (1+C) \ d^q \rceil}^{\infty} (\lambda_{d,s_d,i})^{\varrho} \Big)^{1/\varrho} < \infty.$$

Because of $\lambda_{d,s_d,i} = s_d \lambda_{d,i}$ for any $d,i \in \mathbb{N}$ due to (3.2), this in particular implies that $s_1(\sum_{m=\lceil 1+C \rceil}^{\infty} \lambda_m^{\varrho})^{1/\varrho} \leq C_{\varrho}$ is finite and hence $\lambda = (\lambda_m)_{m \in \mathbb{N}} \in \ell_{\varrho}$. Moreover, we have

$$\sum_{i=1}^{\infty} \left(\lambda_{d,i}\right)^{\varrho} = \|\lambda \mid \ell_{\varrho}\|^{\varrho d}, \quad \text{as well as} \quad \sum_{i=1}^{\lceil (1+C) \, d^q \rceil - 1} \left(\lambda_{d,i}\right)^{\varrho} \leq \lambda_1^{\varrho d} \left(1+C\right) d^q,$$

and therefore

$$\|\lambda \mid \ell_{\varrho}\|^{\varrho d} - \lambda_{1}^{\varrho d} (1+C) d^{q} \leq \left(\frac{C_{\varrho} d^{2q/p}}{s_{d}}\right)^{\varrho} \quad \text{for all } d \in \mathbb{N}.$$
 (3.4)

Now let $\tau > \varrho$ and assume that (III) is violated for this τ . Then $\sup_{d \in \mathbb{N}} s_d \|\lambda \| \ell_{\tau}\|^d$ is infinite since $\lambda \in \ell_{\varrho}$ and $\ell_{\varrho} \hookrightarrow \ell_{\tau}$ with $\|\lambda \| \ell_{\tau}\| < \|\lambda \| \ell_{\varrho}\|$. That means, for any $C_0 \in (0, \infty)$ there necessarily exists a sequence $(d_k)_{k \in \mathbb{N}} \subset \mathbb{N}$ such that for every $k \in \mathbb{N}$,

$$C_0 \le s_{d_k} \|\lambda \mid \ell_{\tau}\|^{d_k} = s_{d_k} \|\lambda \mid \ell_{\varrho}\|^{d_k} / (t_{\varrho,\tau})^{d_k},$$

where we set $t_{\varrho,\tau} = \|\lambda \mid \ell_{\varrho}\|/\|\lambda \mid \ell_{\tau}\| > 1$. Hence, at least for all k larger than a certain $k_0 \in \mathbb{N}$, we conclude that $C_0 \leq s_{d_k} \|\lambda \mid \ell_{\varrho}\|^{d_k}/d_k^{2q/p}$. In particular, we can choose $C_0 > C_{\varrho}$

such that $0 < C_1 = C_{\rho}/C_0 < 1$. Therefore (3.4) implies

$$\|\lambda \| \ell_{\varrho}\|^{\varrho d_k} - \lambda_1^{\varrho d_k} (1+C) d_k^q \le C_1^{\varrho} \|\lambda \| \ell_{\varrho}\|^{\varrho d_k}, \quad k \ge k_0.$$

which leads to

$$\left(1 + \left(\frac{\lambda_2}{\lambda_1}\right)^{\varrho}\right)^{d_k} \le \left(\sum_{m=1}^{\infty} \left(\frac{\lambda_m}{\lambda_1}\right)^{\varrho}\right)^{d_k} = \frac{\|\lambda \mid \ell_{\varrho}\|^{\varrho d_k}}{\lambda_1^{\varrho d_k}} \le C_2 d_k^q$$

for all $k \geq k_0$ and some $C_2 = (1+C)/(1-C_1^{\varrho}) > 0$. Since $\lambda_2 > 0$ and $\varrho > 0$ this is a contradiction and thus we have proven (III) for every $\tau > p/2$. Note that this also shows that

$$\inf\{2\tau \mid \tau \text{ fulfills condition (III)}\} \leq p^*.$$

Step 2. Next we show that (I) follows from (III). Let $\tau > 0$ be given such that (III) holds true and set p = q = r = 0, as well as C = 1. Then, with

$$f(d) = \lceil C(\min\left\{s_d \,\lambda_1^d, 1\right\})^{-p/2} d^q \rceil = 1,$$

we have

$$C_{\tau} = \sup_{d \in \mathbb{N}} \frac{1}{d^{r}} \left(\sum_{i=f(d)}^{\infty} (\lambda_{d,s_{d},i})^{\tau} \right)^{1/\tau} = \sup_{d \in \mathbb{N}} s_{d} \|\lambda \| \ell_{\tau} \|^{d} < \infty.$$

Once more we apply Theorem 2.8 to obtain

$$n_{\text{abs}}^{\text{wor}}(\varepsilon, d; S_{d,s_d}) \leq (1 + C_{\tau}^{\tau}) \varepsilon^{-2\tau}$$
 for all $\varepsilon(0,1]$ and every $d \in \mathbb{N}$.

Thus $S_{(s)}$ is strongly polynomially tractable and

$$p^* \le \inf\{2\tau \mid \tau \text{ fulfills condition (III)}\}.$$

Step 3. The implication (III) \Rightarrow (IV) can be seen as follows. Assume (III) is valid for some $0 < \tau < \infty$ and set $\varrho = \tau$. Then, clearly, $\lambda \in \ell_{\varrho}$. If we now assume (IV) is violated then for any $\delta > 0$ there exists a sequence $(d_k)_{k \in \mathbb{N}} \subset \mathbb{N}$ such that for all k,

$$s_{d_k}^{1/d_k} \ge 1/(\lambda_1 + \delta).$$

Hence, $s_{d_k} \|\lambda \mid \ell_{\tau}\|^{d_k} \ge (\|\lambda \mid \ell_{\tau}\|/(\lambda_1 + \delta))^{d_k}$ tends to infinity (as $k \to \infty$) if we take δ small enough so that $\lambda_1 + \delta < \|\lambda \mid \ell_{\tau}\|$. Since $\tau \in (0, \infty)$ and

$$\lambda_1 < (\lambda_1^{\tau} + \lambda_2^{\tau})^{1/\tau} \le ||\lambda| ||\ell_{\tau}|| < \infty$$

there is some $\alpha_{\tau} \in (0, \infty)$ such that $\|\lambda \mid \ell_{\tau}\| = \lambda_1 + \alpha_{\tau}$. Choosing e.g. $\delta = \alpha_{\tau}/2$ gives the needed contradiction.

Step 4. Finally we have to show that also (IV) \Rightarrow (III). Therefore assume that we have (IV) for some $\varrho \in (0, \infty)$. Then there exist constants $d_0 \in \mathbb{N}$ and $\delta > 0$ such that

$$s_d^{1/d} \le 1/(\lambda_1 + \delta)$$
 for all $d \ge d_0$.

Furthermore note that the function $N(\tau) = \|\lambda \mid \ell_{\tau}\|$ is strictly decreasing and continuous on the interval $[\varrho, \infty]$ and that $N(\varrho) > \lambda_1 = N(\infty)$ because of the ordering of $\lambda = (\lambda_m)_{m \in \mathbb{N}}$. Hence there exists some $\tau \in [\varrho, \infty)$ such that $N(\tau) \leq \lambda_1 + \delta/2$, say. Thus, $\lambda \in \ell_{\tau}$ and for every $d \geq d_0$ we obtain

$$s_d \|\lambda \mid \ell_\tau\|^d \le \left(\frac{\lambda_1 + \delta/2}{\lambda_1 + \delta}\right)^d \le 1.$$

Since the term on the left is also finite for any $d = 1, ..., d_0$ this completes the proof.

Observe that Theorem 3.2 is not very surprising. Indeed, the second assertion in condition (IV) is equivalent to the fact that the dth root of the initial error $\varepsilon_d^{\rm init}$ is asymptotically strictly less than 1. Hence if $S_1: H_1 \to \mathcal{G}_1$ (and thus also the sequence λ) is given then we need to select scaling factors s_d so that $\varepsilon_d^{\rm init} \to 0$, $d \to \infty$, in order to obtain polynomial tractability. More advanced illustrations will be given in Section 3.3.

3.2.2. Weak tractability and the curse. To formulate necessary and sufficient conditions for weak tractability with respect to the worst case setting and the absolute error criterion we need some additional notation. Therefore let $S_{(s)} = (S_{d,s_d})_{d \in \mathbb{N}}$ denote a scaled tensor product problem between Hilbert spaces as explained in Section 3.1 and assume $\lambda_2 > 0$. Then for fixed $d \in \mathbb{N}$ and $0 < \varepsilon < \varepsilon_d^{\text{init}} = s_d^{1/2} \lambda_1^{d/2}$ formula (3.3) implies

$$n(\varepsilon, d) = \min\{n \in \mathbb{N}_0 \mid \lambda_{d, s_d, n+1} \le \varepsilon^2\} = \#\{\boldsymbol{j} \in \mathbb{N}^d \mid s_d \lambda_{j_1} \cdot \ldots \cdot \lambda_{j_d} > \varepsilon^2\}$$
$$= \#\left\{\boldsymbol{j} \in \mathbb{N}^d \mid \frac{\lambda_{j_1}}{\lambda_1} \cdot \ldots \cdot \frac{\lambda_{j_d}}{\lambda_1} > \left(\frac{\varepsilon}{\varepsilon_d^{\text{init}}}\right)^2\right\}. \tag{3.5}$$

By counting the number of indices equal to one we conclude that

$$n(\varepsilon,d) = 1 + \sum_{k=1}^d \binom{d}{k} \# \left\{ \boldsymbol{j} = (j_1,\ldots,j_k) \in (\mathbb{N} \setminus \{1\})^k \; \middle| \; \prod_{l=1}^k \frac{\lambda_{j_l}}{\lambda_1} > \left(\frac{\varepsilon}{\varepsilon_d^{\text{init}}}\right)^2 \right\}.$$

Now we distinguish two cases. First assume that $\lambda_1 = \lambda_2$. Then, obviously, each of the sets in the above equality contains at least one element. Otherwise, in the case $\lambda_1 > \lambda_2$, some of those k-dimensional sets might be empty if k is larger than some $k_d(\varepsilon)$. The reason is that $\lambda_1 > \lambda_2 \ge \lambda_m$ for $m \ge 2$ implies that every factor in $\prod_{l=1}^k \lambda_{j_l}/\lambda_1$ is strictly smaller than 1. In detail, $(\varepsilon/\varepsilon_d^{\rm init})^2 \ge (\lambda_2/\lambda_1)^k$ is equivalent to

$$k > k_d(\varepsilon) = \left\lceil \frac{1}{\ln(\lambda_1/\lambda_2)} \ln\left(\frac{\varepsilon_d^{\text{init}}}{\varepsilon}\right)^2 \right\rceil - 1.$$

Hence, denoting $a_d(\varepsilon) = \min \{d, k_d(\varepsilon)\}$ we have

$$n(\varepsilon, d) = 1 + \sum_{k=1}^{a_d(\varepsilon)} {d \choose k} \# \left\{ \mathbf{j} \in (\mathbb{N} \setminus \{1\})^k \mid \prod_{l=1}^k \frac{\lambda_{j_l}}{\lambda_1} > \left(\frac{\varepsilon}{\varepsilon_d^{\text{init}}}\right)^2 \right\}$$
(3.6)

for $d \in \mathbb{N}$ and $0 < \varepsilon < \varepsilon_d^{\text{init}}$. If $\lambda_1 = \lambda_2$ then the same equality remains true when we formally set $k_d(\varepsilon) = \infty$, i.e. $a_d(\varepsilon) = d$. Moreover, for $d \in \mathbb{N}$ we have $a_d(\varepsilon) = 0$ if and only if $\varepsilon \ge (\lambda_2/\lambda_1)^{1/2} \varepsilon_d^{\text{init}}$. If so, then we obtain $n(\varepsilon, d) = 1$ as long as $\varepsilon < \varepsilon_d^{\text{init}}$ and $n(\varepsilon, d) = 0$ otherwise.

Finally the following statement relates the decay properties of the univariate sequence of the squared singular values λ with the growth behavior of the information complexity $n(\varepsilon,d) = n_{\rm abs}^{\rm wor}(\varepsilon,d;S_{d,s_d})$. It generalizes an assertion given in Novak and Woźniakowski [NW08, p. 178].

LEMMA 3.3. Let $S_{(s)}$ and $\lambda = (\lambda_m)_{m \in \mathbb{N}}$ be as before. Then, for all $\beta \geq 1$,

$$\lambda_n \in o(\ln^{-2\beta} n) \text{ as } n \to \infty \quad \text{if and only if} \quad \ln n(t^{\beta}, 1) \in o(1/t) \text{ as } t \to 0.$$

Proof. Assume $\beta \geq 1$ is fixed and let $t \in (0, (s_1\lambda_2)^{1/(2\beta)})$. Then (3.5) yields, for d = 1,

$$n = n(t^{\beta}, 1) = \min\{n \in \mathbb{N}_0 \mid s_1 \lambda_{n+1} \le t^{2\beta}\} \ge 2.$$

Thus we have $s_1 \lambda_{n(t^{\beta},1)+1} \le t^{2\beta} < s_1 \lambda_{n(t^{\beta},1)}$ and $\ln^{2\beta} n \ge 1/4^{\beta} \cdot \ln^{2\beta} (n+1)$. Combining both these estimates we conclude that

$$\frac{s_1}{4^{\beta}} \frac{\lambda_{n(t^{\beta},1)+1}}{\ln^{-2\beta}(n(t^{\beta},1)+1)} \le \left(\frac{\ln n(t^{\beta},1)}{t^{-1}}\right)^{2\beta} < s_1 \frac{\lambda_{n(t^{\beta},1)}}{\ln^{-2\beta}n(t^{\beta},1)}.$$

Since the one-dimensional information complexity $n(\varepsilon, 1)$ is an increasing function in $1/\varepsilon$, taking the limit for $t \to 0$ proves the claim.

Now we are well-prepared to present necessary conditions for weak tractability based on the representation of the information complexity given in (3.6).

Proposition 3.4. Weak tractability of $S_{(s)}$ implies

$$\lim_{\varepsilon^{-1}+d\to\infty} \frac{\ln \sum_{k=0}^{a_d(\varepsilon)} \binom{d}{k}}{\varepsilon^{-1}+d} = 0 \quad and \quad \lim_{\varepsilon^{-1}+d\to\infty} \frac{\ln n(\varepsilon_1^{\text{init}}\varepsilon/\varepsilon_d^{\text{init}}, 1)}{\varepsilon^{-1}+d} = 0.$$
 (3.7)

If so, then $\lambda_n \in o(\ln^{-2} n)$ as $n \to \infty$. Furthermore,

$$\ln \varepsilon_d^{\text{init}} \in o(d) \quad as \ d \to \infty,$$
 (3.8)

since otherwise $S_{(s)}$ suffers from the curse of dimensionality. If, in addition, $\lambda_1 = \lambda_2$ then we need $\lim_{d\to\infty} \varepsilon_d^{\rm init} = 0$ to avoid the curse. Moreover, in this case weak tractability even yields

$$\varepsilon_d^{\text{init}} \in o(1/d) \quad \text{as } d \to \infty.$$
 (3.9)

Proof. Step 1. We start by proving the necessity of the first limit condition in (3.7) and study its consequences. For this purpose, recall that by the definition of $a_d(\varepsilon)$ all the sets in (3.6) contain at least one element. Consequently, for general $\lambda_1 \geq \lambda_2$ we have

$$n(\varepsilon, d) \ge 1 + \sum_{k=1}^{a_d(\varepsilon)} {d \choose k} = \sum_{k=0}^{a_d(\varepsilon)} {d \choose k} \quad \text{for } d \in \mathbb{N} \text{ and } 0 < \varepsilon < \varepsilon_d^{\text{init}}.$$
 (3.10)

Now assume the existence of a subsequence $(d_l)_{l\in\mathbb{N}}\subset\mathbb{N}$ such that $\varepsilon_{d_l}^{\mathrm{init}}$ grows at least exponentially in d_l for $l\to\infty$. That is, we assume that condition (3.8) is violated. Moreover, fix $\varepsilon=\varepsilon_0\in(0,\inf\{\varepsilon_{d_l}^{\mathrm{init}}\mid l\in\mathbb{N}\})$. Then for any $l\in\mathbb{N}$ and some $\alpha\in(0,1/2)$ the term $a_{d_l}(\varepsilon_0)$ is bounded from below by $\lfloor\alpha\,d_l\rfloor$. Accordingly, (3.10) implies $n(\varepsilon_0,d_l)\geq\binom{d_l}{\lfloor\alpha\,d_l\rfloor}$ for all $l\in\mathbb{N}$. Using similar calculations to those in [NW08, p. 178] we see that this lower bound grows exponentially in d_l . This proves the curse of dimensionality for the scaled problem $S_{(s)}$ and thus contradicts weak tractability.

If we assume in addition that $\lambda_1 = \lambda_2$ then, as already noticed, $a_d(\varepsilon)$ equals d because $k_d(\varepsilon) = \infty$. Thus we obtain $\sum_{k=0}^{a_d(\varepsilon)} \binom{d}{k} = 2^d$ in this case. Therefore the existence of a sequence $(d_l)_{l \in \mathbb{N}}$ such that $\varepsilon_{d_l}^{\text{init}}$ is larger than some C > 0 for all $l \in \mathbb{N}$ would again imply the curse of dimensionality since then we could fix $\varepsilon = \varepsilon_0 = C/2$, say. Moreover $a_d(\varepsilon) = d$ shows that the first part of (3.7) equivalently reads

$$\lim_{\varepsilon^{-1}+d\to\infty} \frac{d}{\varepsilon^{-1}+d} = 0. \tag{3.11}$$

Observe that in any case the term $d/(\varepsilon^{-1}+d)$ is equivalent to the minimum of 1 and εd

(up to some absolute constants). Hence, (3.11) holds true if and only if

$$\lim_{\varepsilon^{-1}+d\to\infty}\varepsilon\,d=0,$$

which in turn is equivalent to $\varepsilon_d^{\text{init}} \in o(1/d)$ for $d \to \infty$. To see this last equivalence, remember that due to Section 1.4 the domain of the sequences $((\varepsilon_k, d_k))_{k \in \mathbb{N}}$ for the limit $\varepsilon^{-1} + d \to \infty$ is restricted per definition to those for which $\varepsilon_k < \varepsilon_{d_k}^{\text{init}}$.

Step 2. We turn to the proof of the second part of (3.7). Again we distinguish the cases $\lambda_1 = \lambda_2$ and $\lambda_1 > \lambda_2$. For the latter case keep in mind that $a_d(\varepsilon) \geq 1$ if and only if $\varepsilon < \varepsilon_d^{\rm init}(\lambda_2/\lambda_1)^{1/2}$. If so, then (3.6) shows that

$$n(\varepsilon, d) \ge 1 + \binom{d}{1} \# \left\{ j \ge 2 \mid \frac{\lambda_j}{\lambda_1} > \left(\frac{\varepsilon}{\varepsilon_d^{\text{init}}}\right)^2 \right\}$$
$$= 1 + d \# \{ j \ge 2 \mid s_1 \lambda_j > (\varepsilon_1^{\text{init}} \varepsilon / \varepsilon_d^{\text{init}})^2 \} \ge n(\varepsilon_1^{\text{init}} \varepsilon / \varepsilon_d^{\text{init}}, 1).$$

On the other hand, if $\varepsilon \in [\varepsilon_d^{\text{init}}(\lambda_2/\lambda_1)^{1/2}, \varepsilon_d^{\text{init}})$ then $n(\varepsilon_1^{\text{init}}\varepsilon/\varepsilon_d^{\text{init}}, 1)$ is no larger than $n(\varepsilon_1^{\text{init}}(\lambda_2/\lambda_1)^{1/2}, 1)$, which is an absolute, positive constant. Thus, as claimed in (3.7), we conclude that

$$0 \leq \frac{\ln n(\varepsilon_1^{\mathrm{init}}\varepsilon/\varepsilon_d^{\mathrm{init}},1)}{\varepsilon^{-1}+d} \leq \max\left\{\frac{\ln n(\varepsilon,d)}{\varepsilon^{-1}+d}, \frac{\ln n(\varepsilon_1^{\mathrm{init}}(\lambda_2/\lambda_1)^{1/2},1)}{\varepsilon^{-1}+d}\right\} \to 0 \tag{3.12}$$

for $\varepsilon^{-1}+d$ tending to infinity in the above sense. In the case $\lambda_1=\lambda_2$ we have $a_d(\varepsilon)=d$, which is trivially bounded from below by 1 for any $\varepsilon\in(0,\varepsilon_d^{\rm init})$. The assertion now follows using the same arguments as in the first part of the previous case. To complete the proof it remains to show that $\lambda_n\in o(\ln^{-2}n)$ as $n\to\infty$. Let us consider the case $d=d_k\equiv 1$ in (3.12). Then we obtain

$$0 \le \frac{\ln n(\varepsilon, 1)}{\varepsilon^{-1}} \le 2 \frac{\ln n(\varepsilon_1^{\text{init}} \varepsilon / \varepsilon_1^{\text{init}}, 1)}{\varepsilon^{-1} + 1} \to 0 \quad \text{as } \varepsilon^{-1} \to \infty.$$

In other words, weak tractability yields $\ln n(\varepsilon, 1) \in o(\varepsilon^{-1})$, which is equivalent to the claimed assertion by Lemma 3.3.

Let us add some comments on the above necessary conditions.

REMARK 3.5. First of all note that from (3.7) we concluded (3.8), which is equivalent to the fact that $\limsup_{d\to\infty} s_d^{1/d} \leq 1/\lambda_1$. Aside from that, (3.7) also implies another condition which we will need later on:

$$\lim_{\varepsilon^{-1}+d\to\infty}\frac{\ln n(\varepsilon_1^{\mathrm{init}}(\varepsilon/\varepsilon_d^{\mathrm{init}})^{1/2},1)}{\varepsilon^{-1}+d}=0.$$

Concluding, we stress that the second part of (3.7) already indicates a certain trade-off between the decay of the sequence $(\lambda_n)_{n\in\mathbb{N}}$ and the growth of the initial error $\varepsilon_d^{\text{init}}$. Indeed, if $(\lambda_n)_{n\in\mathbb{N}}$ decreases almost logarithmically then n(t,1) increases subexponentially as t tends to zero. Consequently, (3.7) can be fulfilled only if $\varepsilon_d^{\text{init}}$ is polynomially bounded in d. On the other hand, if the (squares of the) singular values tend to zero like the inverse of some polynomial, say, then n(t,1) grows polynomially in 1/t and hence it is enough to assume that the initial error is subexponentially bounded in d to fulfill (3.7). \square

We complement the necessary conditions in Proposition 3.4 by the following sufficient conditions for weak tractability of scaled tensor product problems $S_{(s)} = (S_{d,s_d})_{d \in \mathbb{N}}$. For the proof we essentially follow the arguments of Papageorgiou and Petras [PP09] for the unscaled case, which are based on estimates from Woźniakowski [Woź94b].

PROPOSITION 3.6. Let $S_{(s)}$ and $a_d(\varepsilon)$ be as before and assume $\lambda_2 > 0$. If condition (3.7) from Proposition 3.4 holds true and if

$$\lim_{\varepsilon^{-1}+d\to\infty} \frac{a_d(\varepsilon)\ln n(\varepsilon_1^{\text{init}}(\varepsilon/\varepsilon_d^{\text{init}})^{1/2},1)}{\varepsilon^{-1}+d} = 0$$
(3.13)

then $S_{(s)}$ is weakly tractable.

Proof. Given $d \in \mathbb{N}$ and $\varepsilon \in (0, (\lambda_2/\lambda_1)^{1/2} \varepsilon_d^{\text{init}})$ consider the representation (3.6) and keep in mind that for larger ε the information complexity $n(\varepsilon, d)$ is trivially bounded by 1 because then $a_d(\varepsilon) = 0$. For every $k \in \{1, \ldots, a_d(\varepsilon)\}$ we have

$$\# \bigg\{ \boldsymbol{j} \in (\mathbb{N} \setminus \{1\})^k \ \bigg| \ \prod_{l=1}^k \frac{\lambda_{j_l}}{\lambda_1} > \bigg(\frac{\varepsilon}{\varepsilon_d^{\text{init}}} \bigg)^2 \bigg\} \leq \# \bigg\{ \boldsymbol{j} \in \mathbb{N}^{a_d(\varepsilon)} \ \bigg| \ \prod_{l=1}^{a_d(\varepsilon)} \frac{\lambda_{j_l}}{\lambda_1} > \bigg(\frac{\varepsilon}{\varepsilon_d^{\text{init}}} \bigg)^2 \bigg\}$$

since $\lambda_m/\lambda_1 \leq 1$ for all $m \in \mathbb{N}$. Hence we concentrate on all the multi-indices $j = (j_1, \ldots, j_{a_d(\varepsilon)})$ that fulfill

$$\prod_{l=1}^{a_d(\varepsilon)} \frac{\lambda_{j_l}}{\lambda_1} > \left(\frac{\varepsilon}{\varepsilon_d^{\text{init}}}\right)^2. \tag{3.14}$$

Clearly the largest possible index $j_{\max}^{(1)}$ which can appear in those $\mathbf{j} \in \mathbb{N}^{a_d(\varepsilon)}$ is bounded because the sequence $(\lambda_n)_{n \in \mathbb{N}}$ tends to zero as $n \to \infty$. Indeed, using the arguments given in [PP09] we conclude that

$$j_{\max}^{(1)} \le \min \left\{ n \in \mathbb{N}_0 \mid s_1 \lambda_{n+1} \le \left(\varepsilon_1^{\text{init}} \frac{\varepsilon}{\varepsilon_d^{\text{init}}} \right)^2 \right\} = n \left(\varepsilon_1^{\text{init}} \frac{\varepsilon}{\varepsilon_d^{\text{init}}}, 1 \right).$$

More generally, in [PP09] it was noticed that, using the same reasoning, we can bound the *i*th largest index $j_{\max}^{(i)}$ in (3.14) by

$$j_{\text{max}}^{(i)} \le n(\varepsilon_1^{\text{init}}(\varepsilon/\varepsilon_d^{\text{init}})^{1/i}, 1).$$

We use this estimate for i = 1 and 2 to obtain the upper bound

$$a_d(\varepsilon)n(\varepsilon_1^{\mathrm{init}}\varepsilon/\varepsilon_d^{\mathrm{init}},1)n(\varepsilon_1^{\mathrm{init}}(\varepsilon/\varepsilon_d^{\mathrm{init}})^{1/2},1)^{a_d(\varepsilon)-1}$$

for $\#\{j \in \mathbb{N}^{a_d(\varepsilon)} \mid j \text{ fulfills (3.14)}\}$. Note that due to $\varepsilon < \varepsilon_d^{\text{init}}$ both the univariate complexities in the above bound need to be at least 1. Therefore we can extend the estimate by adding an additional factor $n(\varepsilon_1^{\text{init}}(\varepsilon/\varepsilon_d^{\text{init}})^{1/2}, 1)$ and replacing $a_d(\varepsilon)$ by d. In summary we have

$$n(\varepsilon, d) \le dn(\varepsilon_1^{\text{init}} \varepsilon / \varepsilon_d^{\text{init}}, 1) n(\varepsilon_1^{\text{init}} (\varepsilon / \varepsilon_d^{\text{init}})^{1/2}, 1)^{a_d(\varepsilon)} \sum_{k=0}^{a_d(\varepsilon)} \binom{d}{k}$$
(3.15)

for each $d \in \mathbb{N}$ and all $\varepsilon \in (0, (\lambda_2/\lambda_1)^{1/2} \varepsilon_d^{\text{init}})$. Because $n(\varepsilon, d) = 1$ if ε belongs to $[(\lambda_2/\lambda_1)^{1/2} \varepsilon_d^{\text{init}}, \varepsilon_d^{\text{init}})$, the estimate (3.15) remains valid for every $\varepsilon \in (0, \varepsilon_d^{\text{init}})$. Proceeding

as in [PP09] we take the logarithm and divide by $\varepsilon^{-1} + d$ to conclude that

$$\begin{split} \frac{\ln n(\varepsilon,d)}{\varepsilon^{-1}+d} & \leq \frac{\ln d}{\varepsilon^{-1}+d} + \frac{\ln n(\varepsilon_1^{\mathrm{init}}\varepsilon/\varepsilon_d^{\mathrm{init}},1)}{\varepsilon^{-1}+d} \\ & + \frac{a_d(\varepsilon)\ln n(\varepsilon_1^{\mathrm{init}}(\varepsilon/\varepsilon_d^{\mathrm{init}})^{1/2},1)}{\varepsilon^{-1}+d} + \frac{\sum_{k=0}^{a_d(\varepsilon)}\binom{d}{k}}{\varepsilon^{-1}+d}. \end{split}$$

For weak tractability it suffices to show that each of these fractions tends to zero as $\varepsilon^{-1} + d \to \infty$. Obviously, for the first one this is true without any further conditions. For the second and fourth fractions the assertion follows from (3.7). Finally the third fraction tends to zero due to the additional condition (3.13) we imposed for this proposition.

To illustrate our results the following theorem considers several cases of the behavior of the initial error $\varepsilon_d^{\rm init}$.

THEOREM 3.7. Let $S_{(s)} = (S_{d,s_d})_{d \in \mathbb{N}}$ denote a scaled tensor product problem in the sense of Section 3.1. Assume that $\lambda_2 > 0$ and consider the worst case setting with respect to the absolute error criterion.

- Let ln ε_d^{init} ∉ o(d) as d → ∞. Then S_(s) suffers from the curse of dimensionality.
 Let ε_d^{init} ∈ Θ(d^α) as d → ∞ for some α ≥ 0.
- - If $\lambda_1 = \lambda_2$ then $S_{(s)}$ suffers from the curse of dimensionality.
 - In the case $\lambda_1 > \lambda_2$ the problem $S_{(s)}$ is weakly tractable if and only if

$$\lambda_n \in o(\ln^{-2(1+\alpha)} n) \quad \text{as } n \to \infty.$$
 (3.16)

- Let $\varepsilon_d^{\rm init} \to 0$ as $d \to \infty$. Then we never have the curse of dimensionality. Moreover, $S_{(s)}$ is weakly tractable if and only if
 - (i) $\lambda_1 = \lambda_2$ and $\lambda_n \in o(\ln^{-2} n)$ as $n \to \infty$, and $\varepsilon_d^{\rm init} \in o(1/d)$ as $d \to \infty$, or
 - (ii) $\lambda_1 > \lambda_2$ and $\lambda_n \in o(\ln^{-2} n)$ as $n \to \infty$.

Proof. Step 1. In this first step we handle the assertions concerning the curse of dimensionality. From the proof of Proposition 3.4 we know that $S_{(s)}$ suffers from the curse either if $\ln \varepsilon_d^{\text{init}} \notin o(d)$, or if $\lambda_1 = \lambda_2$ and $\lim_{d \to \infty} \varepsilon_d^{\text{init}} \neq 0$. Of course the latter condition is fulfilled particularly if the initial error grows polynomially with the dimension d, i.e. if $\varepsilon_d^{\text{init}} \in \Theta(d^{\alpha})$ for some $\alpha \geq 0$. Furthermore the fact that we cannot have the curse of dimensionality as long as $\varepsilon_d^{\text{init}}$ tends to zero is clear from the definition.

Step 2. Next we show that weak tractability implies (3.16). Note that $\varepsilon_d^{\text{init}} \in \Theta(d^{\alpha})$ implies the existence of some c>0 such that $\varepsilon_d^{\rm init}\geq c\,d^{\alpha}$ for all $d\in\mathbb{N}$. Moreover there is some $d_0 \in \mathbb{N}$ such that $1/c < d^{1+\alpha}$ for every d larger than d_0 . Setting $\varepsilon = 1/d$ now yields

$$\frac{\varepsilon}{\varepsilon_d^{\text{init}}} \le \frac{1}{c} \frac{1}{d^{1+\alpha}} < 1 \quad \text{ for all } d \ge d_0$$

and $\varepsilon^{-1} + d = 2d \to \infty$ as $d \to \infty$. Hence, the sequence $((1/d,d))_{d \ge d_0}$ is admissible for the second limit condition of (3.7). On the other hand,

$$\frac{\ln n(\varepsilon_1^{\mathrm{init}}\varepsilon/\varepsilon_d^{\mathrm{init}},1)}{\varepsilon^{-1}+d} \geq \frac{c'}{2}\,\frac{\ln n((c'/d)^{1+\alpha},1)}{(c'/d)^{-1}} \geq 0$$

where we set $c' = (\varepsilon_1/c)^{1/(1+\alpha)}$. Thus weak tractability implies $\ln n(t^{1+\alpha}, 1) \in o(1/t)$ as $t \to 0$. Now the assertion follows from Lemma 3.3.

Step 3. For polynomial initial errors it remains to prove the converse implication, namely that (3.16) is also sufficient for weak tractability provided that $\lambda_1 > \lambda_2$. To this end, we first show that for all $d \in \mathbb{N}$, every $\varepsilon \in (0, \varepsilon_d^{\text{init}})$, and for some C > 0, we have

$$\varepsilon_d^{\text{init}}/\varepsilon \le C(\varepsilon^{-1} + d)^{1+\alpha}.$$
 (3.17)

To see this, we notice that $\varepsilon_d^{\text{init}} \in \Theta(d^{\alpha})$ implies the existence of some C > 0 such that $\varepsilon_d^{\text{init}} \leq C d^{\alpha}$ for all $d \in \mathbb{N}$. If $\alpha = 0$ then (3.17) is obvious. For $\alpha > 0$ we apply Young's inequality (17) to obtain $\varepsilon_d^{\text{init}}/\varepsilon \leq C(d^{1+\alpha} + (\varepsilon^{-1})^{1+\alpha})$. Now (3.17) follows from the relation $\|\cdot\| \ell_{1+\alpha}\| \leq \|\cdot\| \ell_1\|$, $\alpha \geq 0$, for (two-dimensional) sequence spaces.

We want to deduce weak tractability from Proposition 3.6. Hence we have to check the limit conditions in (3.7) and (3.13). In what follows we abbreviate

$$t = t(\varepsilon, d) = \varepsilon^{-1} + d.$$

Given (3.17), as well as the definition of $a_d(\varepsilon)$ before (3.6), it is easy to see that

$$a_d(\varepsilon) \in \mathcal{O}(\ln t)$$
 as $t \to \infty$.

In particular, $a_d(\varepsilon) < \lfloor \lfloor t \rfloor / 2 \rfloor$ if t is sufficiently large. Moreover, note that d < t implies $\binom{d}{k} \le \binom{\lfloor t \rfloor}{k}$ for all $k \in \{0, 1, \dots, a_d(\varepsilon)\}$ such that $\binom{\lfloor t \rfloor}{a_d(\varepsilon)}$ is an upper bound for each of those binomial coefficients $\binom{d}{k}$. Consequently,

$$\ln \sum_{k=0}^{a_d(\varepsilon)} \binom{d}{k} \le \ln \left((a_d(\varepsilon) + 1) \binom{\lfloor t \rfloor}{a_d(\varepsilon)} \right) \le \ln(2 \, a_d(\varepsilon) \, (e \, \lfloor t \rfloor)^{a_d(\varepsilon)})$$

$$\le \ln 2 + \ln a_d(\varepsilon) + a_d(\varepsilon) (1 + \ln t) \in \mathcal{O}(\ln^2 t) \subseteq o(t)$$

for $t \to \infty$. In other words, the first part of 3.7 is fulfilled. Also the second condition in (3.7) can be shown easily using (3.17). Indeed, by the assumption in (3.16) (or its equivalent reformulation in Lemma 3.3) we conclude that

$$\frac{\ln n(\varepsilon_1^{\mathrm{init}}\varepsilon/\varepsilon_d^{\mathrm{init}},1)}{\varepsilon^{-1}+d} \leq \frac{\ln n(\varepsilon_1^{\mathrm{init}}\,C\,(\varepsilon^{-1}+d)^{-(1+\alpha)},1)}{\varepsilon^{-1}+d} = \frac{1}{C'}\,\frac{\ln n((C'/t)^{1+\alpha},1)}{(C'/t)^{-1}},$$

which tends to zero as $\varepsilon^{-1} + d$ (and therefore also t/C') approaches infinity. Finally,

$$\begin{aligned} a_d(\varepsilon) \ln n(\varepsilon_1^{\text{init}}(\varepsilon/\varepsilon_d^{\text{init}})^{1/2}, 1) &\leq a_d(\varepsilon) \ln n \bigg(\bigg(\frac{C''}{\varepsilon^{-1} + d} \bigg)^{(1+\alpha)/2}, 1 \bigg) \\ &\in \mathcal{O}(\ln t) \mathcal{O}(t^{1/2}) \subseteq o(t) \end{aligned}$$

for $t = \varepsilon^{-1} + d \to \infty$. Hence we have shown (3.13). Now an application of Proposition 3.6 completes the proof for the case of polynomial initial errors.

Step 4. In this last step we consider the case of initial errors which tend to zero for d tending to infinity. We already know from Proposition 3.4 that $\lambda_n \in o(\ln^{-2} n)$ as $n \to \infty$ (or equivalently $\ln n(t, 1) \in o(1/t)$ as $t \to 0$) is necessary for weak tractability,

⁽¹⁷⁾ Recall that Young's inequality states that $a,b \geq 0$ and p,q > 1 with 1/p + 1/q = 1 yield $ab \leq a^p/p + b^q/q$. We use this assertion for $a = d^{\alpha}$, $b = \varepsilon^{-1}$ and $p = 1 + 1/\alpha$, $q = 1 + \alpha$.

independently of the relation of the (squares of the) two largest singular values λ_1 and λ_2 to each other. Moreover Proposition 3.4 states that $\varepsilon_d^{\rm init} \in o(1/d)$ as $d \to \infty$ is a necessary condition when we additionally assume $\lambda_1 = \lambda_2$. It remains to show that these conditions are also sufficient for weak tractability in the particular situations.

If $\lambda_1 > \lambda_2$ then we can exactly follow the lines of Step 3 with $\alpha = 0$ in order to conclude the assertion. Hence we are left with the case $\lambda_1 = \lambda_2$. As in the previous step, we want to apply Proposition 3.6 and thus we need to check the conditions in (3.7) and (3.13). Setting

$$u = u(\varepsilon, d) = \frac{d}{\varepsilon^{-1} + d}$$
(3.18)

we note that (by (3.11) in the proof of Proposition 3.4) u tends to zero as $\varepsilon^{-1} + d \to \infty$. This follows from $\varepsilon_d^{\text{init}} \in o(1/d)$ as $d \to \infty$, and, on the other hand, it implies the first condition in (3.7) because $\sum_{k=0}^{a_d(\varepsilon)} \binom{d}{k}$ equals 2^d . Since, in particular, $\varepsilon_d^{\text{init}} \leq C$ for some C > 0, we have $C' = C/\varepsilon_1^{\text{init}} > 0$ and thus we obtain

$$\frac{\ln n(\varepsilon_1^{\mathrm{init}}\varepsilon/\varepsilon_d^{\mathrm{init}},1)}{\varepsilon^{-1}+d} \leq \frac{\ln n(\varepsilon_1^{\mathrm{init}}/C\varepsilon,1)}{\varepsilon^{-1}+d} \leq C' \frac{\ln n\big((C'(\varepsilon^{-1}+d))^{-1},1\big)}{C'(\varepsilon^{-1}+d)} \to 0$$

as $\varepsilon^{-1} + d \to \infty$, due to $\ln n(t,1) \in o(1/t)$ as $t \to 0$. In other words, we have shown the second condition in (3.7). To see that (3.13) holds true as well, we once more use $\varepsilon_d^{\text{init}} \in o(1/d) \subseteq \mathcal{O}(1/d)$, together with Young's inequality, to conclude that

$$\left(\frac{\varepsilon_d^{\text{init}}}{\varepsilon}\right)^{1/2} \le C_1 \frac{1}{d} d^{1/2} \left(\frac{1}{\varepsilon}\right)^{1/2} \le \frac{C_1}{2} \frac{\varepsilon^{-1} + d}{d}$$

with some $C_1 > 0$. Hence, using (3.18) we have $\varepsilon_1^{\text{init}}(\varepsilon/\varepsilon_d^{\text{init}})^{1/2} \ge C_2 u$ and therefore

$$\frac{d \ln n(\varepsilon_1^{\text{init}}(\varepsilon/\varepsilon_d^{\text{init}})^{1/2}, 1)}{\varepsilon^{-1} + d} \le \frac{1}{C_2} (C_2 u) \ln n(C_2 u, 1) \to 0 \quad \text{as } \varepsilon^{-1} + d \to \infty,$$

because then $C_2 u = C_2 u(\varepsilon, d)$ tends to zero. Since $a_d(\varepsilon) = d$ this yields (3.13) and we deduce weak tractability from Proposition 3.6.

Before we turn to normalized errors we stress that Theorem 3.7 contains at least two surprising results. Firstly, we can have weak tractability even if the initial error of $S_{(s)}$ grows with the dimension. Hence, although the performance of the zero algorithm gets steadily worse for $d \to \infty$ we are not necessarily faced with the curse of dimensionality. In contrast, remember that we need decreasing initial errors in order to obtain polynomial tractability. Secondly, it seems to be quite surprising that also in the case $\lambda_1 = \lambda_2$ we can break the curse by imposing only moderate additional conditions on the scaling sequence s. Indeed, it is enough to guarantee that $\varepsilon_d^{\rm init} = \sqrt{s_d \lambda_1^d} \in o(1/d)$ for $d \to \infty$.

3.2.3. Normalized errors. We complete our studies of the complexity of scaled problems $S_{(s)} = (S_{d,s_d})_{d \in \mathbb{N}}$ by investigating tractability properties with respect to the normalized error criterion. This can be done by analyzing the information complexity of a related problem with respect to absolute errors.

Let $\lambda = (\lambda_n)_{n \in \mathbb{N}}$ and $s = (s_d)_{d \in \mathbb{N}}$ be fixed and define a tensor product problem $T = (T_d: H_d \to \mathcal{G}_d)_{d \in \mathbb{N}}$ from the building blocks $T_1 = ((1/\sqrt{\lambda_1}) S_1): H_1 \to \mathcal{G}_1$ as described in

the proof of Theorem 2.12. Then the extended sequence of squared singular values of T_d , based on the univariate sequence $\mu = (\mu_m)_{m \in \mathbb{N}} = (\lambda_m/\lambda_1)_{m \in \mathbb{N}}$, reads

$$(\mu_{d,1,i})_{i\in\mathbb{N}} = (\mu_{d,i})_{i\in\mathbb{N}} = \left(\frac{\lambda_{d,i}}{\lambda_{d,1}}\right)_{i\in\mathbb{N}} = \left(\frac{\lambda_{d,s_d,i}}{s_d\,\lambda_{d,1}}\right)_{i\in\mathbb{N}} = \left(\frac{\lambda_{d,s_d,i}}{(\varepsilon_d^{\mathrm{init}})^2}\right)_{i\in\mathbb{N}}.$$

Here the second subscript in $\mu_{d,1,i}$ indicates that T can be seen as a trivially scaled tensor product problem. Furthermore, $\varepsilon_d^{\text{init}} = \sqrt{s_d \lambda_1^d}$ denotes the initial error of S_{d,s_d} . Thus, from (3.5) applied to T and $S_{(s)}$ we conclude that

$$n_{\text{abs}}^{\text{wor}}(\varepsilon, d; T_d) = \min \left\{ n \in \mathbb{N}_0 \mid \mu_{d,1,n+1} \le \varepsilon^2 \right\}$$

$$= \min \left\{ n \in \mathbb{N}_0 \mid \lambda_{d,s_d,n+1} \le (\varepsilon \varepsilon_d^{\text{init}})^2 \right\} = n_{\text{abs}}^{\text{wor}}(\varepsilon \varepsilon_d^{\text{init}}, d; S_{d,s_d}).$$
 (3.19)

By definition this also equals $n_{\text{norm}}^{\text{wor}}(\varepsilon, d; S_{d,s_d})$, i.e. the information complexity of S_{d,s_d} with respect to the normalized error criterion. Based on this relation, we can use our results from the previous subsections to prove

THEOREM 3.8. Let $S_{(s)} = (S_{d,s_d})_{d \in \mathbb{N}}$ denote a scaled tensor product problem in the sense of Section 3.1. Assume $\lambda_2 > 0$ and consider the worst case setting with respect to the normalized error criterion.

- Let $\lambda_1 = \lambda_2$. Then $S_{(s)}$ suffers from the curse of dimensionality.
- Let $\lambda_1 > \lambda_2$. Then $S_{(s)}$ is not polynomially tractable. Moreover, in this case $S_{(s)}$ is weakly tractable if and only if $\lambda_n \in o(\ln^{-2} n)$ as $n \to \infty$.

Proof. Note that for all $d \in \mathbb{N}$ the initial error of T_d is 1 since μ_1 , as well as the scaling parameters, equal 1. Thus, obviously, condition (IV) in Theorem 3.2 is violated and therefore T is polynomially intractable with respect to the absolute error criterion. Moreover, the second point of Theorem 3.7 with $\alpha=0$ shows that T suffers from the curse of dimensionality if $\mu_1=\mu_2$. Otherwise, i.e. if $\mu_1>\mu_2$, the problem T is weakly tractable if and only if $\mu_n\in o(\ln^{-2}n)$ as $n\to\infty$. Since we set $\mu_m=\lambda_m/\lambda_1, m\in\mathbb{N}$, all these conditions on $\mu=(\mu_m)_{m\in\mathbb{N}}$ are fulfilled if and only if the corresponding assertions hold true for the sequence $\lambda=(\lambda_m)_{m\in\mathbb{N}}$. Equation (3.19) finally shows that every complexity assertion for T with respect to absolute errors is equivalent to the corresponding statement for $S_{(s)}$ and the normalized error criterion. This simple observation completes the proof. \blacksquare

In conclusion, the scaling sequence $s = (s_d)_{d \in \mathbb{N}}$ does not have any influence on the complexity of $S_{(s)}$, as long as we consider normalized errors. So the advantages of scaling are completely ruled out in this setting.

3.3. Examples. In this last part of Chapter 3 we briefly discuss two applications of the complexity results obtained in the previous section. We start by proving that our assertions reproduce the known facts for unscaled tensor product problems studied in Theorems 2.11 and 2.12.

EXAMPLE 3.9 (Unscaled problems). Let $S_{(s)} = (S_{d,s_d})_{d \in \mathbb{N}}$ denote a tensor product problem between Hilbert spaces in the sense of Section 3.1 where all the scaling factors s_d equal 1. As usual we assume $\lambda_2 > 0$ and consider the worst case setting. Then for every $d \in \mathbb{N}$ the operators S_{d,s_d} coincide with S_d as defined in Section 2.4.1. Since we already

saw that for the normalized error criterion the conditions stated in Theorem 3.8 exactly match the assertions of Theorem 2.12, it remains to consider the absolute error criterion. Here $\varepsilon_d^{\text{init}}$ is given by $\lambda_1^{d/2}$. Hence there are three scenarios for the behavior of the initial error depending on the largest squared singular value λ_1 of the underlying operator S_1 .

From Theorem 3.2 we know that strong polynomial tractability and polynomial tractability are equivalent; see (I) and (II), respectively. Moreover, condition (IV) shows that this holds if and only if $\lambda_1 < 1$ and $\lambda = (\lambda_m)_{m \in \mathbb{N}} \in \ell_{\tau}$ for some $\tau \in (0, \infty)$. In this case the exponent of strong polynomial tractability is given by

$$p^* = \inf \left\{ 2\tau \mid \sup_{d \in \mathbb{N}} \|\lambda \mid \ell_\tau\|^d < \infty \right\} = \inf \left\{ 2\tau \mid \left(\sum_{m=1}^\infty (\lambda_m)^\tau \right)^{1/\tau} \le 1 \right\}.$$

In turn, $\lambda_1 \geq 1$ yields polynomial intractability. More precisely, if $\lambda_1 > 1$ then the initial error grows exponentially in d and $S = S_{(s)}$ suffers from the curse of dimensionality due to the first point of Theorem 3.7. Setting $\alpha = 0$ the second point of the latter theorem describes the case of constant initial errors, which corresponds to the case $\lambda_1 = 1$ in the unscaled situation. In detail, if $\lambda_2 = \lambda_1 = 1$ then we are faced with the curse again. In contrast, if $\lambda_2 < 1$ then we have weak tractability if and only if

$$\lambda_n \in o(\ln^{-2} n) \quad \text{as } n \to \infty.$$
 (3.20)

Finally the initial error tends to zero exponentially fast if $\lambda_1 < 1$. The last point of Theorem 3.7 thus shows that in this case the curse of dimensionality is not possible and that (3.20) is necessary and sufficient for weak tractability.

Altogether these results exactly match the conditions stated in Theorem 2.11. Hence, scaled tensor product problems indeed yield a generalization. \Box

So let us turn to a more advanced application. To this end, recall the definition of $S_{d,s_d} \colon \mathcal{F}_d \to \mathcal{G}_d$ in Section 3.1. There we constructed the source spaces \mathcal{F}_d by scaling the norm in the tensor product space $H_d = H_1 \otimes \cdots \otimes H_1$. Alternatively we can think of \mathcal{F}_d as the successively taken tensor product of some building blocks $H^{(k)}$, $k = 1, \ldots, d$, in the sense of Section 2.4.1, where we define $H^{(k)}$ to be the univariate space H_1 scaled by some factor $s^{(k)} > 0$. That is, let

$$\langle \cdot, \cdot \rangle_{H^{(k)}} = \frac{1}{s^{(k)}} \, \langle \cdot, \cdot \rangle_{H_1}.$$

Then the scaling factor s_d in dimension d is given by $\prod_{k=1}^d s^{(k)} > 0$. The following example illustrates how the behavior of the generator sequence $(s^{(k)})_{k \in \mathbb{N}}$ affects the complexity of $S_{(s)}$.

EXAMPLE 3.10. Because scaling has no influence on assertions for normalized errors we restrict ourselves to the absolute error criterion in what follows. For simplicity we further assume that $\lambda_1 > \lambda_2 > 0$ and that the generator sequence is non-increasing, i.e.

$$s^{(1)} \ge \dots \ge s^{(k)} \ge \dots > 0, \quad k \in \mathbb{N}.$$

Then Theorem 3.2 states that $S_{(s)}$ is strongly polynomially tractable if and only if the geometric mean of the first d elements $s^{(k)}$ is asymptotically strictly smaller than $1/\lambda_1$, provided that $\lambda \in \ell_{\tau}$ for some $\tau > 0$. This holds iff at most finitely many of these

generators are bounded from below by $1/\lambda_1$. Moreover, from Remark 3.5 we know that we need

$$\limsup_{d \to \infty} s_d^{1/d} = \limsup_{d \to \infty} \Big(\prod_{k=1}^d s^{(k)}\Big)^{1/d} \leq \frac{1}{\lambda_1}$$

in order to obtain weak tractability. Therefore let the generators be given by

$$s^{(k)} = \frac{1}{\lambda_1} (1 + \delta_k), \quad k \in \mathbb{N},$$

with a non-increasing null sequence $(\delta_k)_{k\in\mathbb{N}}$ and note that then some elementary calculations yield

$$\exp\left(\frac{c}{2}\sum_{k=1}^{d}\delta_{k}\right) \leq \varepsilon_{d}^{\text{init}} \leq \exp\left(\frac{1}{2}\sum_{k=1}^{d}\delta_{k}\right), \quad d \in \mathbb{N},$$

where $c = \ln(1 + \delta_1)/\delta_1 \le 1$. Furthermore this observation shows that $\varepsilon_d^{\text{init}} \in \Theta(d^{\alpha})$ as $d \to \infty$, for some $\alpha \ge 0$, implies that

$$L = \lim_{d \to \infty} \frac{1}{\ln d} \sum_{k=1}^{d} \delta_k \in 2 \alpha \left[1, \frac{1}{c} \right].$$

Conversely, from the existence of L it follows that for any $\delta > 0$ there is some $d_0 = d_0(\delta)$ such that

$$\varepsilon_d^{\text{init}} \in [d^{\alpha_1}, d^{\alpha_2}] \quad \text{ for all } d \ge d_0,$$

where $\alpha_1 = cL/2 - \delta$ and $\alpha_2 = L/2 + \delta$. Hence, if L is sufficiently small then the initial error $\varepsilon_d^{\text{init}}$ behaves like a polynomial of small degree and thus a quite slow decay of the sequence $(\lambda_n)_{n \in \mathbb{N}}$ is enough to derive weak tractability using Theorem 3.7. \square

4. Problems on function spaces with weighted norms

In [NW09] it is shown that the approximation problem defined on $C^{\infty}([0,1]^d)$ is intractable. In fact, Novak and Woźniakowski considered the linear space F_d of all real-valued infinitely differentiable functions f defined on the unit cube $[0,1]^d$ in d dimensions for which the norm

$$||f| F_d|| = \sup_{\alpha \in \mathbb{N}_0^d} ||D^{\alpha} f| L_{\infty}([0, 1]^d)||$$
(4.1)

of $f \in F_d$ is finite. In this case the (uniform) approximation problem is given by the sequence of solution operators $S = (S_d)_{d \in \mathbb{N}}$,

$$S_d = \mathrm{id}_d \colon \widetilde{F}_d \to L_\infty([0,1]^d), \quad f \mapsto \mathrm{id}_d(f) = f, \quad d \in \mathbb{N},$$
 (4.2)

defined on the unit ball $\widetilde{F}_d = \mathcal{B}(F_d)$ of F_d . The authors studied this problem in the worst case setting using algorithms from the classes $\mathcal{A}_d^{n,\text{cont}}$ and $\mathcal{A}_d^{n,\text{adapt}}$ as defined in Section 1.3.

The *initial error* of this problem is given by $\varepsilon_d^{\text{init}} = e^{\text{wor}}(0, d; \text{id}_d) = 1$, the norm of the embedding $F_d \hookrightarrow L_{\infty}$, since $A_{0,d} \equiv 0$ is a valid choice of an algorithm which does not use any information on f; see Proposition 2.4. This means that the problem is well-scaled such that there is no difference in studying the absolute or the normalized error criterion.

Now [NW09, Theorem 1] implies that the nth minimal worst case error of L_{∞} -approximation defined on F_d satisfies

$$e^{\text{wor}}(n, d; \text{id}_d) = 1$$
 for all $n = 0, 1, \dots, 2^{\lfloor d/2 \rfloor} - 1$. (4.3)

Therefore, for all $d \in \mathbb{N}$ and every $\varepsilon \in (0,1)$, the information complexity is bounded from below by

$$n^{\text{wor}}(\varepsilon, d; \text{id}_d) > 2^{\lfloor d/2 \rfloor}$$
.

Hence the problem suffers from the curse of dimensionality; in particular it is intractable. One possibility of avoiding this exponential dependence on d, i.e. to break the curse, is to shrink the function space F_d by introducing weights.

In the present chapter we follow this idea. We show that turning to spaces equipped with product weights can dramatically improve the tractability behavior of certain problems such as uniform approximation. In Section 4.1 we formally introduce the concept of weighted spaces by considering the examples of weighted Banach spaces of smooth functions and of weighted reproducing kernel Hilbert spaces. Upper error bounds for uniform approximation in the latter class of spaces are then studied in Section 4.2. Afterwards, in Section 4.3, we show how to use those bounds for the L_{∞} -approximation problem defined on scales of smooth functions. Moreover we prove corresponding lower bounds on the

information complexity which enable us to give necessary and sufficient conditions for several kinds of tractability in terms of the used weights. Most of the results stated in this chapter are published in [Wei12b].

4.1. The concept of weighted spaces. The idea to introduce weights directly into the norm of the function space appeared for the first time in a paper of Sloan and Woźniakowski in 1998; see [SW98]. They studied the integration problem defined over some Sobolev Hilbert space, equipped with so-called *product weights*, to explain the overwhelming success of QMC integration rules. Thenceforth weighted problems attracted a lot of attention.

For example it turned out that tractability of approximation of linear compact operators between Hilbert spaces can be fully characterized in terms of the weights and the singular values of the operators if we use information operations from the class $\Lambda^{\rm all}$. The proof of this kind of assertions is once again based on the singular value decomposition; see Section 2.3. One such result is given in Section 4.4 below.

But first let us illustrate the concept of weighted spaces by modifying the space F_d we introduced before.

4.1.1. Weighted Banach spaces of smooth functions. A closer look at the norm given in (4.1) shows that for $f \in \mathcal{B}(F_d)$ we have

$$||D^{\alpha}f| L_{\infty}([0,1]^d)|| \le 1 \quad \text{for all } \alpha \in \mathbb{N}_0^d.$$

$$(4.4)$$

Hence every derivative is equally important. In order to shrink the space, for each $\alpha \in \mathbb{N}_0^d$ we replace the right-hand side of inequality (4.4) by a non-negative weight γ_{α} . For α with $|\alpha| = 1$ this means that we control the importance of every single variable. So, the norm in the weighted space F_d^{γ} is now given by

$$||f| F_d^{\gamma}|| = \sup_{\alpha \in \mathbb{N}_0} \frac{1}{\gamma_{\alpha}} ||D^{\alpha} f| L_{\infty}([0, 1]^d)||, \tag{4.5}$$

where we demand $D^{\alpha}f$ to be equal to zero if $\gamma_{\alpha} = 0$. It is clear from the construction that we indeed shrink the space if all γ_{α} are chosen strictly less than one.

Since this approach is quite general we restrict ourselves to so-called *product weights* (with uniformly bounded generators) in what follows. Thus we assume that for every $d \in \mathbb{N}$ there exists an ordered and uniformly bounded sequence

$$C_{\gamma} \ge \gamma_{d,1} \ge \cdots \ge \gamma_{d,d} \ge 0.$$

Then for $d \in \mathbb{N}$ the product weight sequence $\gamma = (\gamma_{\alpha})_{\alpha \in \mathbb{N}_0^d}$ is given by

$$\gamma_{\alpha} = \prod_{j=1}^{d} (\gamma_{d,j})^{\alpha_j}, \quad \alpha \in \mathbb{N}_0^d.$$
 (4.6)

Note that the influence of x_j on f is now controlled by the so-called generator weight $\gamma_{d,j}$. Since $\gamma_{d,j}=0$ for some $j\in\{1,\ldots,d\}$ implies that f does not depend on x_j,\ldots,x_d we assume that $\gamma_{d,d}>0$ for the rest of this chapter. Moreover observe that the ordering of $\gamma_{d,j}$ does not restrict generality. Later on we will see that tractability of our problem will only depend on summability properties of the generator weights.

Among other things, we show in Section 4.3.3 that for the L_{\infty}-approximation problem defined on the Banach spaces F_d^{γ} with the norm given above and generator weights $\gamma_{d,j} \equiv \gamma^{(j)} \in \Theta(j^{-\beta})$ we have

- intractability for $\beta = 0$,
- weak tractability but no polynomial tractability for $0 < \beta < 1$,
- strong polynomial tractability for $1 < \beta$.

Furthermore, we prove that for $\beta = 1$ the problem is not strongly polynomially tractable.

4.1.2. Weighted Hilbert spaces and weighted RKHS. Let us briefly discuss the idea of weighted norms in the case of Hilbert (function) spaces, before we turn to weighted RKHSs. Our approach is based on a generalization of the so-called ANOVA (18) decomposition of d-variate functions f, where d is an arbitrarily large integer. For the ease of presentation we follow the lines of [NW08, Section 5.3.1]. Thus, we focus our attention on Hilbert function spaces constructed from tensor products and equipped with some assumptions that can be significantly relaxed. For further information on more general settings the interested reader is referred to [KSWW10b] and the references therein.

Given a d-fold tensor product space $H_d = H_1 \otimes \cdots \otimes H_1$, $d \in \mathbb{N}$, as well as an orthonormal basis $\{e_i \mid i \in \mathbb{N}\}$ of the underlying univariate Hilbert space (19) H_1 that contains the constant function $e_1 \equiv 1$, it is easy to see that every $f \in H_d$ can be represented as

$$f = \sum_{\mathfrak{u} \subseteq \{1, \dots, d\}} f_{\mathfrak{u}}.$$

In this decomposition the (formally d-variate) functions $f_{\mathfrak{u}}$ solely depend on the variables x_j with index $j \in \mathfrak{u}$. The main advantage of this kind of representation is that for fixed f the collection of all $f_{\mathfrak{u}}$, $\mathfrak{u} \subseteq \{1,\ldots,d\}$, can be taken mutually orthogonal with respect to the inner product $\langle\cdot,\cdot\rangle_{H_d}$ in H_d . Therefore the norm of $f \in H_d$ can be expressed by

$$||f| H_d||^2 = \sum_{\mathfrak{u} \subseteq \{1, \dots, d\}} ||f_{\mathfrak{u}}| H_d||^2 = \sum_{\mathfrak{u} \subseteq \{1, \dots, d\}} ||f_{\mathfrak{u}, 1}| H_{|\mathfrak{u}|}||^2,$$

where $f_{\mathfrak{u},1}$ equals $f_{\mathfrak{u}}$ interpreted as an element of the $|\mathfrak{u}|$ -fold tensor product space $H_{|\mathfrak{u}|}$ of the closed subspace

$$H_1' = \{ h \in H_1 \mid \langle h, e_1 \rangle_{H_1} = 0 \} \subset H_1$$

with itself. That is, in the unweighted situation the contribution of each $f_{\mathfrak{u}}$ to the norm of $f \in H_d$ is the same.

Now suppose that we have some additional a priori knowledge about the importance of some (groups of) variables in dimension d. This can be modeled by assigning positive (20)

⁽¹⁸⁾ analysis of variance.

 $^(^{19})$ We assume H_1 to be separable and infinite-dimensional in order to keep the notation as short as possible.

^{(&}lt;sup>20</sup>) Also zero weights are possible but for reasons of simplification we do not discuss this more complicated situation in the present brief introduction to weighted Hilbert spaces.

weights $\gamma_{d,\mathfrak{u}}$ to each of the 2^d subsets \mathfrak{u} of $\{1,\ldots,d\}$. We denote the collection of these weights by $\gamma_{(d)} = \{\gamma_{d,\mathfrak{u}} \mid \mathfrak{u} \subseteq \{1,\ldots,d\}\}$. Then it can be verified that

$$\langle f, g \rangle_{\gamma_{(d)}} = \sum_{\mathfrak{u} \subset \{1, \dots, d\}} \frac{1}{\gamma_{d, \mathfrak{u}}} \langle f_{\mathfrak{u}}, g_{\mathfrak{u}} \rangle_{H_d} = \sum_{\mathfrak{u} \subset \{1, \dots, d\}} \frac{1}{\gamma_{d, \mathfrak{u}}} \langle f_{\mathfrak{u}, 1}, g_{\mathfrak{u}, 1} \rangle_{H_{|\mathfrak{u}|}}$$
(4.7)

defines an inner product on the tensor product space H_d which implies an equivalent norm depending on $\gamma_{(d)}$. The Hilbert space H_d furnished with this new inner product will be denoted by $H_d^{\gamma_{(d)}}$. At this point we need to stress the fact that for general weights $\gamma_{(d)}$ these spaces are no longer tensor product spaces, although their construction is based on $H_d = H_1 \otimes \cdots \otimes H_1$ and $H_{|\mathfrak{u}|}$, respectively. To overcome this problem we restrict ourselves to the case of product weights in the following. Thus we assume

$$\gamma_{d,\mathfrak{u}} = \prod_{k \in \mathfrak{u}} \gamma_{d,k} \tag{4.8}$$

for some positive $\gamma_{d,k}$, $k=1,\ldots,d$, and every $\mathfrak{u}\subseteq\{1,\ldots,d\}$. Then it can be checked that indeed $H_d^{\gamma_{(d)}}$ is again a tensor product space. For a study of other types of weights such as *finite-order*, *finite-diameter*, *order-dependent* or the recently developed POD (²¹) weights we refer to Novak and Woźniakowski [NW08, Section 5.3.2] and to Kuo, Schwab and Sloan [KSS11].

In the last decade it turned out that weighted norms provide a powerful tool to vanquish the curse of dimensionality that we are often faced with. Since the $H_d^{\gamma_{(d)}}$'s are still Hilbert spaces the complexity analysis of the weighted problems $S^{\gamma_{(d)}} = (S_d^{\gamma_{(d)}})$: $\mathcal{B}(H_d^{\gamma_{(d)}}) \to \mathcal{G}_d)_{d \in \mathbb{N}}$ is again based on the singular value decomposition presented in Section 2.3.1; at least in the cases where the target spaces \mathcal{G}_d are also Hilbert spaces. Fortunately, the introduced weights enter the spectrum of the operator $W_d^{\gamma_{(d)}} = (S_d^{\gamma_{(d)}})^{\dagger} S_d^{\gamma_{(d)}}$ in a straightforward way. Therefore in many cases tractability properties of S can be fully characterized in terms of the singular values and the introduced weights.

For our purposes weighted Hilbert spaces that possess a reproducing kernel are of particular interest. Typical examples of such weighted RKHSs are the following unanchored Sobolev spaces endowed with product weights which will play an important role in our further argumentation; see also Sloan and Woźniakowski [SW02]. Instead of applying the presented approach which is based on decompositions we use the common procedure and define them directly.

EXAMPLE 4.1 (Unanchored Sobolev spaces \mathcal{H}_d^{γ}). As usual we start with the definition for d=1 and $\gamma>0$. Then the space \mathcal{H}_1^{γ} is nothing but the Sobolev space of all absolutely continuous real-valued functions f defined on the unit interval [0,1] whose first derivative (2^2) f' belongs to the space $L_2([0,1])$. The difference from the classical Sobolev space is the inner product, which here depends on the parameter γ :

$$\langle f, g \rangle_{\mathcal{H}_{1}^{\gamma}} = \langle f, g \rangle_{L_{2}([0,1])} + \gamma^{-1} \langle f', g' \rangle_{L_{2}([0,1])}$$

$$= \int_{0}^{1} f(x) g(x) d\lambda^{1}(x) + \gamma^{-1} \int_{0}^{1} f'(x) g'(x) d\lambda^{1}(x), \quad f, g \in \mathcal{H}_{1}^{\gamma}.$$
(4.9)

^{(&}lt;sup>21</sup>) **p**roduct and **o**rder-**d**ependent.

^{(&}lt;sup>22</sup>) In the weak or distributional sense.

For the sake of completeness we define the space \mathcal{H}_1^0 as the limit of \mathcal{H}_1^{γ} for $\gamma \to 0$. Consequently, the derivatives of $f \in \mathcal{H}_1^0$ need to vanish λ^1 -almost everywhere on [0,1], which implies that the space \mathcal{H}_1^0 only consists of constant functions. This coincides with the common convention 0/0 = 0.

Note that the univariate space \mathcal{H}_1^{γ} algebraically coincides with its anchored analogue $\widetilde{\mathcal{H}}_1^{\gamma}$ where the term $\langle f,g \rangle_{\mathbf{L}_2([0,1])}$ in (4.9) is replaced by f(a)g(a) for some anchor point $a \in [0,1]$. For details we refer to [SW02] and [Wei12b]. Finally we mention that for positive parameters γ all these definitions imply equivalent norms on the classical Sobolev space $W_2^1([0,1])$.

Once more the d-variate spaces \mathcal{H}_d^{γ} for d>1 are defined by a tensor product construction similar to that in Section 2.4.1. We set $\mathcal{H}_d^{\gamma} = \bigotimes_{k=1}^d \mathcal{H}_1^{\gamma_{d,k}}$, where now γ denotes a (subset of a) product weight sequence $(\gamma_{\alpha})_{\alpha \in \{0,1\}^d}$ induced by some generator weights $\gamma_{d,k}$, $k=1,\ldots,d$; see (4.6). Remember that at the beginning of this chapter we assumed $\gamma_{d,d}>0$ for all $d\in\mathbb{N}$. That is, we avoid taking the trivial spaces \mathcal{H}_1^0 as factors in the definition of \mathcal{H}_d^{γ} .

How does the inner product of \mathcal{H}_d^{γ} look like? Following the lines of Section 2.4.1 it is uniquely determined by the coordinatewise inner products of the factors of simple tensors $f = \bigotimes_{k=1}^d f_k$ and $g = \bigotimes_{k=1}^d g_k$, where $f_k, g_k \in \mathcal{H}_1^{\gamma_{d,k}}$ for $k = 1, \ldots, d$. Consequently,

$$\begin{split} \langle f,g\rangle_{\mathcal{H}_{d}^{\gamma}} &= \prod_{k=1}^{d} \langle f_{k},g_{k}\rangle_{\mathcal{H}_{1}^{\gamma_{d,k}}} = \prod_{k=1}^{d} \left(\langle f_{k},g_{k}\rangle_{\mathbf{L}_{2}([0,1])} + \frac{1}{\gamma_{d,k}} \langle f'_{k},g'_{k}\rangle_{\mathbf{L}_{2}([0,1])}\right) \\ &= \sum_{\mathfrak{u}\subseteq\{1,\ldots,d\}} \prod_{k\in\mathfrak{u}} \frac{1}{\gamma_{d,k}} \prod_{k\in\mathfrak{u}} \langle f'_{k},g'_{k}\rangle_{\mathbf{L}_{2}([0,1])} \prod_{j\in\{1,\ldots,d\}\setminus\mathfrak{u}} \langle f_{j},g_{j}\rangle_{\mathbf{L}_{2}([0,1])} \\ &= \sum_{\mathfrak{u}\subseteq\{1,\ldots,d\}} \prod_{k\in\mathfrak{u}} \frac{1}{\gamma_{d,k}} \int_{[0,1]^{d}} \prod_{k\in\mathfrak{u}} f'_{k}(x_{k}) g'_{k}(x_{k}) \prod_{j\in\{1,\ldots,d\}\setminus\mathfrak{u}} f_{j}(x_{j}) g_{j}(x_{j}) \, \mathrm{d}\lambda^{d}(\boldsymbol{x}) \\ &= \sum_{\mathfrak{u}\in\{1,\ldots,d\}} \frac{1}{\gamma_{d,\mathfrak{u}}} \int_{[0,1]^{d}} \frac{\partial^{|\mathfrak{u}|} f}{\partial x_{\mathfrak{u}}}(\boldsymbol{x}) \frac{\partial^{|\mathfrak{u}|} g}{\partial x_{\mathfrak{u}}}(\boldsymbol{x}) \, \mathrm{d}\lambda^{d}(\boldsymbol{x}), \end{split}$$

where we used (4.8) and the shorthand notation $\partial^{|\mathfrak{u}|}/\partial x_{\mathfrak{u}}$ for $\prod_{k\in\mathfrak{u}}\partial/\partial x_k$. Note that this representation resembles (4.7) from the general approach to weighted Hilbert spaces introduced at the beginning of this subsection. For our purposes it is more convenient to rewrite the subsets $\mathfrak{u}\subseteq\{1,\ldots,d\}$ in terms of multi-indices $\boldsymbol{\alpha}=(\alpha_1,\ldots,\alpha_d)\in\{0,1\}^d$. In detail, we set $\alpha_k=1$ if $k\in\mathfrak{u}$ and $\alpha_k=0$ otherwise. Then we can express the norm of any $f\in\mathcal{H}_d^{\gamma}$ by

$$||f| \mathcal{H}_d^{\gamma}||^2 = \sum_{\alpha \in \{0,1\}^d} \frac{1}{\gamma_{\alpha}} \int_{[0,1]^d} |D^{\alpha} f(\boldsymbol{x})|^2 d\lambda^d(\boldsymbol{x})$$
(4.10)

since then $\gamma_{\mathfrak{u}} = \gamma_{\alpha}$. The inner products of the multivariate anchored spaces, \mathcal{H}_d^{γ} , can be found by a similar reasoning; see [Wei12b, p. 67] for the final result.

It is known (cf. Micchelli and Wahba [MW81]) that the univariate spaces \mathcal{H}_1^{γ} are reproducing kernel Hilbert spaces for any $\gamma > 0$. Consequently, this property is transferred

to the multivariate tensor product space. To stress this fact we write $\mathcal{H}(K_d^{\gamma})$ for \mathcal{H}_d^{γ} in what follows. Equation (5) in [WW09] now states that the reproducing kernel

$$K_d^{\gamma} \colon [0,1]^d \times [0,1]^d \to \mathbb{R}$$

in dimension $d \ge 1$ is given by $(^{23})$

$$K_d^{\gamma}(\boldsymbol{x}, \boldsymbol{y}) = \prod_{k=1}^d \frac{\sqrt{\gamma_{d,k}}}{\sinh(\sqrt{\gamma_{d,k}})} \cosh(\sqrt{\gamma_{d,k}} \left(1 - \max\left\{x_k, y_k\right\}\right)) \cosh(\sqrt{\gamma_{d,k}} \min\left\{x_k, y_k\right\}),$$

 $x, y \in [0, 1]^d$. For d = 1 this kernel formula follows from Thomas–Agnan [Tho96, Corollary 2] whereas the higher-dimensional generalization for product weights γ results from the tensor product structure; see (2.20) in Section 2.5. In particular we note that K_d^{γ} is continuous (and thus also bounded) along its diagonal

$$\{(\boldsymbol{x}, \boldsymbol{y}) \in [0, 1]^{2d} \mid \boldsymbol{x} = \boldsymbol{y}\}.$$

Moreover, from [WW09, Lemma 4.1] we know that for $\gamma > 0$ the set

$$E_1(\gamma) = \{e_{1,\gamma,i} \colon [0,1] \to \mathbb{R} \mid i \in \mathbb{N}\}\$$

with $e_{1,\gamma,1} \equiv 1$ and

$$e_{1,\gamma,i}(x) = \cos(\pi(i-1)x)\sqrt{\frac{2\gamma}{\gamma + \pi^2(i-1)^2}}, \quad x \in [0,1], i \ge 2,$$

is an orthonormal basis in the univariate space $\mathcal{H}(K_1^{\gamma})$. Applying the arguments from Section 2.4.1 this leads to an ONB $E_d(\gamma)$ of $\mathcal{H}(K_d^{\gamma}) = \bigotimes_{k=1}^d \mathcal{H}(K_1^{\gamma_{d,k}})$ that consists of the tensor product functions

$$\widetilde{e}_{d,\gamma,\boldsymbol{m}} = \bigotimes_{k=1}^{d} e_{1,\gamma_{d,k},m_k}, \quad \boldsymbol{m} = (m_1,\dots,m_d) \in \mathbb{N}^d.$$
 (4.11)

For a direct proof of this result we refer to [NW08, Appendix A.2.1] (²⁴) and to [WW09, Lemma 4.2]. Actually, these proofs show a little more, namely that the functions $\tilde{e}_{d,\gamma,\boldsymbol{m}}$ together with

$$\widetilde{\lambda}_{d,\gamma,\boldsymbol{m}} = \prod_{k=1}^{d} \lambda_{1,\gamma_{d,k},m_k} = \prod_{k=1}^{d} \frac{\gamma_{d,k}}{\gamma_{d,k} + \pi^2(m_k - 1)^2}, \quad \boldsymbol{m} \in \mathbb{N}^d,$$
(4.12)

describe the full set of eigenpairs $\{(\widetilde{\lambda}_{d,\gamma,\boldsymbol{m}},\widetilde{e}_{d,\gamma,\boldsymbol{m}}) \mid \boldsymbol{m} \in \mathbb{N}^d\}$ of the operator $W_d^{\gamma} = (S_d^{\gamma})^{\dagger} S_d^{\gamma}$ where $S_d^{\gamma} \colon \mathcal{H}_d^{\gamma} \hookrightarrow L_2([0,1]^d)$ denotes the solution operator of the L₂-approximation problem on $\mathcal{H}_d^{\gamma} = \mathcal{H}(K_d^{\gamma})$. \square

4.2. Uniform approximation in reproducing kernel Hilbert spaces. The main result of this section is based on a paper of Kuo, Wasilkowski and Woźniakowski [KWW08]. In contrast to the presentation given in [Wei12b] we decided to apply this result to the case of unanchored Sobolev space introduced in Section 4.1.2 instead of the anchored analogue studied in [KWW08]. This opens up the opportunity to explain the underlying

⁽²³⁾ Here sinh and cosh denote the hyperbolic sine and cosine functions, respectively.

^{(&}lt;sup>24</sup>) Note the missing factor 1/2 in [NW08, p. 351, line 5].

ideas without literally repeating the proof given in [KWW08] while obtaining a result which is (according to our knowledge) not published elsewhere so far.

We start with an upper error bound which remains valid for any reproducing kernel Hilbert space $\mathcal{H}(K_d)$ of real-valued functions f on $[0,1]^d$ with

$$\operatorname{ess\,sup}_{\boldsymbol{x}\in[0,1]^d} K_d(\boldsymbol{x},\boldsymbol{x}) < \infty. \tag{4.13}$$

This condition guarantees that $\mathcal{H}(K_d)$ is continuously embedded into $L_{\infty}([0,1]^d)$ since the reproducing property (2.17), together with the Hahn–Banach theorem (cf. [Yos80, IV.6, Cor.2]), implies that $\|\mathrm{id}_d |\mathcal{L}(\mathcal{H}(K_d), L_{\infty}([0,1]^d))\|$ is given by

$$\sup_{f \in \mathcal{B}(\mathcal{H}(K_d))} \|f \mid \mathcal{L}_{\infty}([0,1]^d)\| = \underset{\boldsymbol{x} \in [0,1]^d}{\operatorname{ess \, sup}} \sup_{f \in \mathcal{B}(\mathcal{H}(K_d))} |f(\boldsymbol{x})|$$

$$= \underset{\boldsymbol{x} \in [0,1]^d}{\operatorname{ess \, sup}} \sup_{f \in \mathcal{B}(\mathcal{H}(K_d))} |\langle f, K_d(\cdot, \boldsymbol{x}) \rangle_{\mathcal{H}(K_d)}|$$

$$= \underset{\boldsymbol{x} \in [0,1]^d}{\operatorname{ess \, sup}} K_d(\boldsymbol{x}, \boldsymbol{x})^{1/2}.$$

Now the upper bound reads as follows:

PROPOSITION 4.2. For $d \in \mathbb{N}$ consider a RKHS $\mathcal{H}(K_d)$, where K_d fulfills (4.13), i.e. $\mathcal{H}(K_d) \hookrightarrow L_{\infty}([0,1]^d)$. Furthermore, suppose $\Xi = \{\xi_j : [0,1]^d \to \mathbb{R} \mid j \in \mathbb{N}\}$ is some orthonormal basis of $\mathcal{H}(K_d)$ and let $n \in \mathbb{N}_0$. Then the algorithm $A_{n,d}^\Xi \in \mathcal{A}_d^{n,\text{lin}}(\Lambda^{\text{all}})$ given by

$$f \mapsto A_{n,d}^{\Xi} f = \sum_{j=1}^{n} \langle f, \xi_j \rangle_{\mathcal{H}(K_d)} \, \xi_j(\cdot)$$

for uniform approximation on $\mathcal{H}(K_d)$ fulfills

$$\Delta^{\text{wor}}\left(A_{n,d}^{\Xi}; \text{id}_d \colon \mathcal{B}(\mathcal{H}(K_d)) \to \mathcal{L}_{\infty}([0,1]^d)\right) \le \left\| \sum_{j=n+1}^{\infty} \xi_j(\cdot)^2 \mid \mathcal{L}_{\infty}([0,1]^d) \right\|^{1/2}. \tag{4.14}$$

Proof. Since Ξ builds an ONB we may represent any $f \in \mathcal{H}(K_d)$ by its basis expansion, $f = \sum_{j=1}^{\infty} \langle f, \xi_j \rangle_{\mathcal{H}(K_d)} \xi_j$. Therefore Parseval's identity implies

$$|f(\boldsymbol{x}) - A_{n,d}^{\Xi} f(\boldsymbol{x})| = |(f - A_{n,d}^{\Xi} f)(\boldsymbol{x})| = \left| \sum_{j=n+1}^{\infty} \langle f, \xi_j \rangle_{\mathcal{H}(K_d)} \xi_j(\boldsymbol{x}) \right|$$
$$= \left| \left\langle f, \sum_{j=n+1}^{\infty} \xi_j(\boldsymbol{x}) \xi_j \right\rangle_{\mathcal{H}(K_d)} \right|,$$

which can be estimated from above using the Cauchy-Schwarz inequality. Thus

$$\left| f(\boldsymbol{x}) - A_{n,d}^{\Xi} f(\boldsymbol{x}) \right| \leq \| f \mid \mathcal{H}(K_d) \| \left\| \sum_{j=n+1}^{\infty} \xi_j(\boldsymbol{x}) \, \xi_j \mid \mathcal{H}(K_d) \right\|$$

$$= \| f \mid \mathcal{H}(K_d) \| \left(\sum_{j=n+1}^{\infty} \xi_j(\boldsymbol{x})^2 \right)^{1/2}$$

$$(4.15)$$

for every $f \in \mathcal{H}(K_d)$ and all fixed $x \in [0,1]^d$. Taking the (essential) supremum with respect to x in the d-dimensional unit cube and the supremum over all $f \in \mathcal{B}(\mathcal{H}(K_d))$ gives the desired result. \blacksquare

We note in passing that we can easily prove more than we stated in the last assertion. In what follows we only need the upper error bound given above, so that we restrict ourselves to some brief comments on further results in the next remark.

REMARK 4.3. For fixed $\boldsymbol{x} \in [0,1]^d$ we see that the function $f^* = C \sum_{j=n+1}^{\infty} \xi_j(\boldsymbol{x}) \, \xi_j$ with C > 0 gives equality in (4.15). Of course, we can choose the constant C such that $\|f^* \mid \mathcal{H}(K_d)\| = 1$ provided that \boldsymbol{x} is not a common root of ξ_j for all j > n. Hence, the upper bound in (4.14) is sharp.

Moreover, [KWW08, Theorem 2] shows that the *n*th minimal worst case error for L_{∞} -approximation on $\mathcal{H}(K_d)$ is given by

$$e^{\text{wor}}(n, d; \text{id}_d : \mathcal{B}(\mathcal{H}(K_d)) \to L_{\infty}([0, 1]^d)) = \inf_{\Xi = \{\xi_j | j \in \mathbb{N}\}} \left\| \sum_{j=n+1}^{\infty} \xi_j(\cdot)^2 \mid L_{\infty}([0, 1]^d) \right\|^{1/2},$$

where the infimum is taken with respect to all orthonormal bases $\Xi \subset \mathcal{H}(K_d)$. Thus, any clever choice of the basis Ξ in Proposition 4.2 leads to algorithms $A_{n,d}^{\Xi}$ with almost optimal worst case errors. \Box

Next we apply Proposition 4.2 to the weighted unanchored Sobolev spaces \mathcal{H}_d^{γ} introduced in Section 4.1.2 using the basis $\Xi = E_d(\gamma)$ given in (4.11). Since the ordering of the basis functions $\xi \in \Xi$ is essential for our application we rearrange them non-increasingly with respect to their L_{∞} -norm:

$$\|\xi_i \mid \mathcal{L}_{\infty}([0,1]^d)\| \ge \|\xi_{i+1} \mid \mathcal{L}_{\infty}([0,1]^d)\| \quad \text{for all } j \in \mathbb{N}.$$
 (4.16)

We obtain an estimate which resembles the corresponding result for the anchored case studied in [Wei12b, Proposition 2].

COROLLARY 4.4. For $n \in \mathbb{N}_0$ and $d \in \mathbb{N}$ there exists an algorithm $A_{n,d}^* \in \mathcal{A}_d^{n,\text{lin}}(\Lambda^{\text{all}})$ for uniform approximation on \mathcal{H}_d^{γ} such that for every $\tau \in (1/2,1)$,

$$\Delta^{\mathrm{wor}}(A_{n,d}^*; \mathrm{id}_d \colon \mathcal{B}(\mathcal{H}_d^{\gamma}) \to \mathrm{L}_{\infty}([0,1]^d)) < a_{\tau} \exp\left(b_{\tau} \sum_{k=1}^d (\gamma_{d,k})^{\tau}\right) n^{-(1-\tau)/(2\tau)},$$

where the constants $a_{\tau}, b_{\tau} > 0$ are independent of γ , n, and d.

Proof. To keep the notation as short as possible we abbreviate the L_{∞} -norm in d dimensions, $\|\cdot| L_{\infty}([0,1]^d)\|$, by $\|\cdot\|_d$ within this proof.

Following our plan we fix $n \in \mathbb{N}_0$, as well as $d \in \mathbb{N}$, and take $A_{n,d}^* = A_{n,d}^\Xi$ defined in Proposition 4.2 with $\Xi = E_d(\gamma)$ as above. From (4.11) we conclude for d = 1 and any $\gamma > 0$ that

$$||e_{1,\gamma,1}^2||_1 = 1$$
 and $||e_{1,\gamma,i}^2||_1 = \frac{2\gamma}{\gamma + \pi^2(i-1)^2} < \frac{2\gamma}{\pi^2}(i-1)^{-2}, \quad i \ge 2.$

Moreover, for every simple tensor $f = \bigotimes_{k=1}^d f_k \in \mathcal{H}(K_d^{\gamma})$ we clearly have

$$||f||_d = \prod_{k=1}^d ||f_k||_1$$
 and $f(\boldsymbol{x})^2 = \prod_{k=1}^d f_k(x_k)^2$, $\boldsymbol{x} \in [0, 1]^d$.

Consequently, for any $j \in \mathbb{N}$ and all $\tau \in (1/2, \infty)$ the ordering of Ξ given in (4.16) implies

$$\begin{split} j\|\xi_{j}^{2}\|_{d}^{\tau} &\leq \sum_{m=1}^{\infty} \|\xi_{m}^{2}\|_{d}^{\tau} = \sum_{\boldsymbol{m} \in \mathbb{N}^{d}} \|\widetilde{e}_{d,\gamma,\boldsymbol{m}}^{2}\|_{d}^{\tau} = \prod_{k=1}^{d} \sum_{i=1}^{\infty} \|e_{1,\gamma_{d,k},i}^{2}\|_{1}^{\tau} \\ &= \prod_{k=1}^{d} \left(1 + \sum_{i=2}^{\infty} \|e_{1,\gamma_{d,k},i}^{2}\|_{1}^{\tau}\right) < \prod_{k=1}^{d} \left(1 + \left(\frac{2\gamma_{d,k}}{\pi^{2}}\right)^{\tau} \sum_{i=2}^{\infty} (i-1)^{-2\tau}\right) \\ &= \prod_{k=1}^{d} \left(1 + c_{\tau}\gamma_{d,k}^{\tau}\right), \end{split}$$

where we set $c_{\tau} = (2/\pi^2)^{\tau} \zeta(2\tau)$. Hence, if $\tau \in (1/2, 1)$ then

$$\left\| \sum_{j=n+1}^{\infty} \xi_j^2 \right\|_d \le \sum_{j=n+1}^{\infty} \|\xi_j^2\|_d < \sum_{j=n+1}^{\infty} j^{-1/\tau} \left(\prod_{k=1}^d \left(1 + c_\tau \gamma_{d,k}^\tau \right) \right)^{1/\tau} < \infty.$$

Since the first factor is no larger than $\int_n^\infty x^{-1/\tau} d\lambda^1(x) = (\tau/(1-\tau))n^{-(1-\tau)/\tau}$ and the second factor can be bounded by $\exp((c_\tau/\tau) \cdot \sum_{k=1}^d (\gamma_{d,k})^\tau)$ we conclude that

$$\left\| \sum_{j=n+1}^{\infty} \xi_j(\cdot)^2 \mid \mathcal{L}_{\infty}([0,1]^d) \right\|^{1/2} < a_{\tau} \exp\left(b_{\tau} \sum_{k=1}^d (\gamma_{d,k})^{\tau}\right) n^{-(1-\tau)/(2\tau)}$$

with $a_{\tau} = \sqrt{\tau/(1-\tau)}$ and $b_{\tau} = c_{\tau}/(2\tau) = (2/\pi^2)^{\tau} \zeta(2\tau)/(2\tau)$. Now the claim follows from (4.14) in Proposition 4.2.

- **4.3.** Uniform approximation in Banach spaces of smooth functions. Our derivation of necessary and sufficient conditions for various kinds of tractability for the L_{∞} -approximation problem defined on the weighted spaces F_d^{γ} introduced in Section 4.1.1 is based on simple embedding arguments. To present them, we consider a whole scale of Banach spaces \mathcal{F}_d^{γ} (of which F_d^{γ} is a special case). Then we first study lower bounds on the *n*th minimal error on a space $\mathcal{P}_d^{\gamma} \hookrightarrow \mathcal{F}_d^{\gamma}$ which consists of *d*-variate polynomials of low degree. Afterwards, in Section 4.3.2, we use the results for $\mathcal{H}_d^{\gamma} \hookleftarrow \mathcal{F}_d^{\gamma}$ from Section 4.2 to deduce corresponding upper bounds. Finally we discuss a couple of concrete examples in Section 4.3.3.
- **4.3.1. Lower bounds for spaces of low-degree polynomials.** Following the lines of [Wei12b, Section 4] we use Proposition 2.2 to obtain a lower bound for the L_{∞} -approximation error for the space

$$\mathcal{P}_{d}^{\gamma} = \operatorname{span}\left\{p_{i} : [0, 1]^{d} \to \mathbb{R}, p_{i}(\boldsymbol{x}) = \boldsymbol{x}^{i} = \prod_{i=1}^{d} (x_{j})^{i_{j}} \mid i = (i_{1}, \dots, i_{d}) \in \{0, 1\}^{d}\right\}$$

of all real-valued d-variate polynomials of degree at most one in each coordinate direction, defined on the unit cube $[0,1]^d$. We equip this linear space with the weighted norm

$$||f| \mathcal{P}_{d}^{\gamma}|| = \max_{\alpha \in \{0,1\}^{d}} \frac{1}{\gamma_{\alpha}} ||D^{\alpha}f| L_{\infty}([0,1]^{d})||, \quad f \in \mathcal{P}_{d}^{\gamma},$$
(4.17)

similar to (4.5), where γ is a product weight sequence as described in (4.6), and study the worst case setting.

THEOREM 4.5. For $d \in \mathbb{N}$ and $n \in \mathbb{N}_0$ assume $A_{n,d} \in \mathcal{A}_d^{n,\text{cont}} \cup \mathcal{A}_d^{n,\text{adapt}}$ to be an arbitrary algorithm for the uniform approximation problem defined on \mathcal{P}_d^{γ} . Then

$$\Delta^{\text{wor}}(A_{n,d}; \text{id}_d : B_r(\mathcal{P}_d^{\gamma}) \to L_{\infty}([0,1]^d)) \ge r \quad \text{for all } r \ge 0$$

provided that $n < 2^s$, where $s = s(\gamma, d) \in \{0, 1, ..., d\}$ is some integer such that

$$s > \frac{1}{2 + C_{\gamma}} \left(\sum_{j=1}^{d} \gamma_{d,j} - 2 \right).$$
 (4.18)

Proof. The proof of this lower error bound consists of several steps. First we fix $d \in \mathbb{N}$ and construct a partition of the set of coordinates $\{1, \ldots, d\}$ into s+1 parts which we will need later and with $s = s(\gamma, d)$ satisfying (4.18). In a second step we define a special linear subspace $V \subseteq \mathcal{P}_d^{\gamma}$ with dim $V = 2^s$. Step 3 then shows that V satisfies the assumptions of Proposition 2.2. The proof is completed in Step 4.

Step 1. For $k \in \{0, ..., d\}$ let us define inductively $m_0 = 0$ and

$$m_k = \inf \left\{ t \in \mathbb{N} \mid m_{k-1} < t \le d, \text{ with } 2 \le \sum_{j=m_{k-1}+1}^t \gamma_{d,j} \right\}$$

with the usual convention inf $\emptyset = \infty$. Note that the infimum coincides with the minimum in the finite case, since then $m_k \in \mathbb{N}$. Moreover we set

$$s = \max \{k \in \{0, \dots, d\} \mid m_k < \infty\}.$$

We let $I_k = \{m_{k-1} + 1, m_{k-1} + 2, \dots, m_k\}$ for $k = 1, \dots, s$. This gives a unique disjoint partition of the set

$$\{1,\ldots,d\} = \left(\bigcup_{k=1}^{s} I_{k}\right) \cup \{m_{s}+1,\ldots,d\},$$

and m_k denotes the last element of the block I_k . For all $k = 1, \ldots, s$ we conclude that

$$2 \le \sum_{j \in I_k} \gamma_{d,j} < 2 + \gamma_{d,m_k} \le 2 + C_{\gamma},$$

where C_{γ} is the uniform upper bound for $\gamma_{d,j}$; see Section 4.1.1. Finally, summation of those inequalities gives

$$\sum_{j=1}^{d} \gamma_{d,j} < \sum_{k=1}^{s} \sum_{j \in I_k} \gamma_{d,j} + 2 < (2 + C_{\gamma})s + 2,$$

and (4.18) follows immediately.

If s=0 then we can stop at this point since the initial error is 1 as the norm of the embedding $\mathcal{P}_d^{\gamma} \hookrightarrow L_{\infty}$ (cf. Proposition 2.4) and the remaining assertion is trivial. Hence, from now on we can assume that s>0 and thus $m_s\geq 1$.

Step 2. To apply Proposition 2.2 we have to construct a linear subspace V of $\mathcal{F} = \mathcal{P}_d^{\gamma}$ such that condition (2.1) holds for the target space $\mathcal{G} = L_{\infty}([0,1]^d)$, the embedding operator $S = \mathrm{id}_d$, and a = 1. Subsequently, we restrict ourselves to the subset

$$\widehat{\mathcal{F}} = \{ f \in \mathcal{F} \mid f \text{ depends only on } x_1, \dots, x_{m_s} \}$$

of \mathcal{F} , since we can interpret $\widehat{\mathcal{F}}$ as the space $\mathcal{P}_{m_s}^{\gamma}$ by a simple isometric isomorphism.

We are ready to construct a suitable space V using the partition from Step 1. We define V as the span of all functions $g_i \colon [0,1]^{m_s} \to \mathbb{R}$, $i = (i_1, \ldots, i_s) \in \{0,1\}^s$, of the form

$$g_{\boldsymbol{i}}(\boldsymbol{x}) = \prod_{k=1}^{s} \left(\sum_{j \in I_k} \gamma_{d,j} x_j \right)^{i_k}, \quad \boldsymbol{x} \in X = [0,1]^{m_s}.$$

Clearly, V is a linear subspace of $\mathcal{P}_{m_s}^{\gamma}$ and with the interpretation above it is also a linear subspace of \mathcal{F} . Moreover it is easy to see that we have by construction

$$||g| \mathcal{F}|| = ||g| \mathcal{P}_{m_0}^{\gamma}||$$
 and $||g| L_{\infty}(X)|| = ||g| L_{\infty}([0,1]^d)||$ for $g \in V$.

Finally we note that dim $V = \#\{0,1\}^s = 2^s$. It remains to show that this subspace is the right choice to prove the claim using Proposition 2.2.

Step 3. The proof of condition (2.1) needed, i.e.,

$$||g| \mathcal{P}_{m_{-}}^{\gamma}|| \leq ||g| \mathcal{L}_{\infty}(X)||$$
 for all $g \in V$,

is a little bit technical. By the special structure of the functions $g \in V$, the left-hand side reduces to $\max\{\gamma_{\alpha}^{-1} \| D^{\alpha}g \mid \mathcal{L}_{\infty}(X) \| \mid \alpha \in \mathbb{M}\}$, where the maximum is taken over all multi-indices α in the set

$$\mathbb{M} = \Big\{ \boldsymbol{\alpha} \in \{0, 1\}^{m_s} \ \Big| \ \sum_{j \in I_k} \alpha_j \le 1 \text{ for all } k = 1, \dots, s \Big\}.$$

This is simply because for $\alpha \notin \mathbb{M}$ we have $D^{\alpha}g \equiv 0$, and then the inequality is trivial. To simplify the notation let us define

$$T: \{0,1\}^{m_s} \to \mathbb{N}_0^s, \quad \boldsymbol{\alpha} \mapsto T(\boldsymbol{\alpha}) = \boldsymbol{\sigma} = (\sigma_1, \dots, \sigma_s),$$

where

$$\sigma_k = \sum_{j \in I_k} \alpha_j$$
 for $k = 1, \dots, s$.

Note that $T(\mathbb{M}) = \{0,1\}^s$. Moreover, for every $g = \sum_{i \in \{0,1\}^s} c_i g_i(\cdot) \in V$ we define a function

$$h_g \colon Z = \underset{k=1}{\overset{s}{\times}} \left[0, \sum_{j \in I_k} \gamma_{d,j} \right] \to \mathbb{R}, \quad \boldsymbol{z} \mapsto h_g(\boldsymbol{z}) = \sum_{\boldsymbol{i} \in \{0,1\}^s} c_{\boldsymbol{i}} \prod_{k=1}^s z_k^{i_k} = \sum_{\boldsymbol{i} \in \{0,1\}^s} c_{\boldsymbol{i}} \, \boldsymbol{z}^{\boldsymbol{i}}.$$

Hence, $h_g(z) = g(x)$ under the transformation $x \mapsto z$ such that

$$z_k = \sum_{j \in I_k} \gamma_{d,j} x_j$$
 for every $k = 1, \dots, s$ and every $\boldsymbol{x} \in X$.

The span, W, of all functions $h: Z \to \mathbb{R}$ with this structure is a linear space, too. Furthermore, elementary calculations yield

$$(D_{\boldsymbol{x}}^{\boldsymbol{\alpha}}g)(\boldsymbol{x}) = \left(\prod_{j=1}^{m_s} (\gamma_{d,j})^{\alpha_j}\right) (D_{\boldsymbol{z}}^{T(\boldsymbol{\alpha})} h_g)(\boldsymbol{z})$$
(4.19)

for all $g \in V$, $\alpha \in \mathbb{M}$ and $x \in X$. Here the x and z in D_x^{α} and $D_z^{T(\alpha)}$ indicate differentiation with respect to x and z, respectively. Since the mapping $x \mapsto z$ is surjective we obtain $\|D^{\alpha}g \mid \mathcal{L}_{\infty}(X)\| = \gamma_{\alpha}\|D^{T(\alpha)}h_g \mid \mathcal{L}_{\infty}(Z)\|$ by the form of γ given in (4.6). Thus,

$$\max_{\alpha \in \mathbb{M}} \frac{1}{\gamma_{\alpha}} \|D^{\alpha} g \mid \mathcal{L}_{\infty}(X)\| = \max_{\sigma \in \{0,1\}^{s}} \|D^{\sigma} h_{g} \mid \mathcal{L}_{\infty}(Z)\|.$$

Observe that (4.19) with $\alpha = 0$ yields in particular $||g| | L_{\infty}(X)|| = ||h_g| | L_{\infty}(Z)||$. Therefore the claim reduces to

$$\max_{\sigma \in \{0,1\}^s} \|D^{\sigma} h_g \mid \mathcal{L}_{\infty}(Z)\| \le \|h_g \mid \mathcal{L}_{\infty}(Z)\| \quad \text{ for every } g \in V.$$

We show this estimate for every $h \in W$, i.e.,

$$||D^{\boldsymbol{\sigma}}h| | \mathcal{L}_{\infty}(Z)|| \le ||h| | \mathcal{L}_{\infty}(Z)|| \quad \text{for all } \boldsymbol{\sigma} \in \{0, 1\}^s.$$

$$(4.20)$$

We start with the special case of one derivative. That is, we first consider $\sigma = e_k$ for a certain $k \in \{1, ..., s\}$. Since h is affine in each coordinate we can represent it as

$$h(\boldsymbol{z}) = a(\boldsymbol{z}_{(\boldsymbol{k})})z_k + b(\boldsymbol{z}_{(\boldsymbol{k})})$$

with functions a and b which only depend on $z_{(k)} = (z_1, \ldots, z_{k-1}, z_{k+1}, \ldots, z_s)$. Hence we have $(D^{e_k}h)(z) = a(z_{(k)})$ and we need to show that

$$|a(\boldsymbol{z}_{(\boldsymbol{k})})| \le \max \left\{ |b(\boldsymbol{z}_{(\boldsymbol{k})})|, \left| a(\boldsymbol{z}_{(\boldsymbol{k})}) \sum_{j \in I_k} \gamma_{d,j} + b(\boldsymbol{z}_{(\boldsymbol{k})}) \right| \right\}. \tag{4.21}$$

Obviously this is true for every $z \in Z$ with $a(z_{(k)}) = 0$. For $a(z_{(k)}) \neq 0$ we can divide by $|a(z_{(k)})|$ to get

$$1 \le \max \left\{ \left| t \right|, \left| \sum_{j \in I_k} \gamma_{d,j} - t \right| \right\}$$

if we set $t = -b(z_{(k)})/a(z_{(k)})$. The last maximum is minimal if both of its entries coincide. This is for $t = \frac{1}{2} \sum_{j \in I_k} \gamma_{d,j}$. Consequently, we need to ensure that

$$2 \le \sum_{j \in I_k} \gamma_{d,j}$$

to obtain (4.21) for all admissible $z \in Z$. But this is true for every $k \in \{1, ..., s\}$ by definition of the sets I_k in Step 1. Thus we have shown (4.20) for the special case $\sigma = e_k$ for all $k \in \{1, ..., s\}$.

Inequality (4.20) also holds true for every $\sigma \in \{0,1\}^s$ by an easy inductive argument on the cardinality of $|\sigma|$. Indeed, if $|\sigma| \geq 2$ then $\sigma = \sigma' + e_k$ with $|\sigma'| = |\sigma| - 1$. We now need to estimate $||D^{\sigma'+e_k}h|| L_{\infty}(Z)||$. Since $(D^{e_k}h)(z) = a(z_{(k)})$ has the same structure as the function h itself, we see that $||D^{\sigma'+e_k}h|| L_{\infty}(Z)||$ equals $||D^{\sigma'}a(z_{(k)})|| L_{\infty}(Z)||$ and the proof of (4.20) is then completed by the inductive step.

Step 4. Collecting the previous equalities and estimates we obtain

$$||g| \mathcal{P}_{d}^{\gamma}|| = ||g| \mathcal{P}_{m_{s}}^{\gamma}|| = \max_{\substack{\boldsymbol{\alpha} \in \{0,1\}^{m_{s}} \\ T(\boldsymbol{\alpha}) \in \{0,1\}^{s}}} \frac{1}{\gamma_{\boldsymbol{\alpha}}} ||D^{\boldsymbol{\alpha}}g| L_{\infty}(X)|| = \max_{\boldsymbol{\sigma} \in \{0,1\}^{s}} ||D^{\boldsymbol{\sigma}}h_{g}| L_{\infty}(Z)||$$

$$\leq ||h_{g}| L_{\infty}(Z)|| = ||g| L_{\infty}(X)|| = ||g| L_{\infty}([0,1]^{d})||$$

for every $g \in V$, where V is a linear subspace of $\mathcal{F} = \mathcal{P}_d^{\gamma}$ with $\dim V = 2^s$. Therefore Proposition 2.2 with a = 1 yields that for $n < \dim V$ the worst case error

$$\Delta^{\mathrm{wor}}(A_{n,d}; \mathrm{id}_d \colon B_r(\mathcal{P}_d^{\gamma}) \to \mathrm{L}_{\infty}([0,1]^d))$$

of any algorithm $A_{n,d}$ from the class $\mathcal{A}_d^{n,\mathrm{cont}} \cup \mathcal{A}_d^{n,\mathrm{adapt}}$ is lower bounded by r, the radius of the centered ball $B_r(\mathcal{P}_d^{\gamma})$.

4.3.2. Complexity results via embeddings. Keeping in mind the assertions shown in the previous sections, we are ready to give conditions for tractability of the uniform approximation problem

$$\mathrm{App} = (\mathrm{App}_d)_{d \in \mathbb{N}}, \quad \mathrm{App}_d \colon \mathcal{B}(\mathcal{F}_d^{\gamma}) \to \mathrm{L}_{\infty}([0,1]^d), \quad \mathrm{App}_d(f) = \mathrm{id}_d(f) = f.$$

We suppose that $(\mathcal{F}_d^{\gamma})_{d\in\mathbb{N}}$ is a sequence of Banach spaces of real-valued functions f defined on the unit cube $[0,1]^d$. We further assume that this sequence depends on product weights $\gamma = (\gamma_{\alpha})_{\alpha \in \mathbb{N}_0^d}$ and fulfills one of the following simple assumptions:

(A4.1)
$$\mathcal{P}_d^{\gamma} \hookrightarrow \mathcal{F}_d^{\gamma}$$
 with norm

$$C_{1,d} \le cd^{q_1}$$
 for all $d \in \mathbb{N}$

and some absolute constants $c, q_1 \geq 0$,

(A4.2) $\mathcal{F}_d^{\gamma} \hookrightarrow \mathcal{H}_d^{\gamma}$ with norm

$$C_{2,d} \le a \exp\left(b \sum_{j=1}^{d} (\gamma_{d,j})^t\right) \quad \text{for all } d \in \mathbb{N}$$
 (4.22)

and some absolute constants a > 0, $b \ge 0$, as well as a parameter $t \in (0,1]$ independent of d and γ .

Here the spaces \mathcal{P}_d^{γ} and $\mathcal{H}_d^{\gamma} = \mathcal{H}(K_d^{\gamma})$ are defined as in Sections 4.1.2 and 4.3.1, respectively.

To simplify the notation we use the commonly known definitions of the so-called sum exponents (25) for the product weight sequence $\gamma = (\gamma_{\alpha})_{\alpha \in \mathbb{N}_0^d}$, $d \in \mathbb{N}$, induced by uniformly bounded generator weights $0 < \gamma_{d,j} \le C_{\gamma}$, $j = 1, \ldots, d$; see (4.6). We set

$$p(\gamma) = \inf \left\{ \kappa \ge 0 \mid P_{\kappa}(\gamma) = \limsup_{d \to \infty} \sum_{j=1}^{d} (\gamma_{d,j})^{\kappa} < \infty \right\},$$
$$q(\gamma) = \inf \left\{ \kappa \ge 0 \mid Q_{\kappa}(\gamma) = \limsup_{d \to \infty} \frac{\sum_{j=1}^{d} (\gamma_{d,j})^{\kappa}}{\ln(d+1)} < \infty \right\},$$

with the usual convention that $\inf \emptyset = \infty$.

 $^(^{25})$ Note that some authors use the name decay for $1/p(\cdot).$

The following necessary conditions for (strong) polynomial tractability slightly generalize Theorem 2 of [Wei12b].

PROPOSITION 4.6 (Necessary conditions). Assume that (A4.1) holds true for some non-negative q_1 . Consider L_{∞} -approximation over $(\mathcal{F}_d^{\gamma})_{d \in \mathbb{N}}$ in the worst case setting with respect to the class of algorithms $\mathcal{A}_d^{n,\mathrm{cont}} \cup \mathcal{A}_d^{n,\mathrm{adapt}}$ and the absolute error criterion. Then

$$n^{\text{wor}}(\varepsilon, d; \text{App}_d) > \frac{1}{2} \cdot 2^{\wedge} \left(\frac{1}{2 + C_{\gamma}} \sum_{i=1}^{d} \gamma_{d,i} \right)$$
 (4.23)

for all $d \in \mathbb{N}$ and every $\varepsilon \in (0, C_{1,d}^{-1})$. Hence

- if the problem App is polynomially tractable then $q(\gamma) \leq 1$;
- if $q_1 = 0$ and the problem is strongly polynomially tractable then $p(\gamma) \leq 1$.

Proof. Let $d \in \mathbb{N}$. Due to (A4.1), every algorithm $A_{n,d} \in \mathcal{A}_d^{n,\text{cont}} \cup \mathcal{A}_d^{n,\text{adapt}}$ for L_{∞} -approximation defined on \mathcal{F}_d^{γ} also applies to the embedded space \mathcal{P}_d^{γ} . Furthermore the embedding constant $C_{1,d}$ implies that the ball $B_r(\mathcal{P}_d^{\gamma})$ of radius $r = C_{1,d}^{-1}$ in \mathcal{P}_d^{γ} is completely contained in the unit ball $\mathcal{B}(\mathcal{F}_d^{\gamma})$ of \mathcal{F}_d^{γ} . Therefore,

$$\Delta^{\text{wor}}(A_{n,d}; \operatorname{App}_d: \mathcal{B}(\mathcal{F}_d^{\gamma}) \to \operatorname{L}_{\infty}([0,1]^d)) \ge \Delta^{\text{wor}}(A_{n,d}|_{\mathcal{P}^{\gamma}}; \operatorname{id}_d: B_r(\mathcal{P}_d^{\gamma}) \to \operatorname{L}_{\infty}([0,1]^d)).$$

From Theorem 4.5 we see that the latter quantity is lower bounded by $r = C_{1,d}^{-1}$ provided that $n < 2^s$, where $s = s(\gamma, d) \in \{0, \ldots, d\}$ satisfies (4.18). Since this lower bound holds for any such $A_{n,d}$ it remains valid for the *n*th minimal error, i.e.

$$e^{\text{wor}}(n, d; \text{App}_d) \ge C_{1,d}^{-1}$$
 for all $n < 2^s$.

Hence we obtain $n^{\mathrm{wor}}(\varepsilon, d; \mathrm{App}_d) \geq 2^s$ for all $d \in \mathbb{N}$ and every $\varepsilon \in (0, C_{1,d}^{-1})$, which implies (4.23) using (4.18).

Now suppose the problem App = $(App_d)_{d\in\mathbb{N}}$ is polynomially tractable. Then there are constants C, p > 0 and $q_2 \geq 0$ such that

$$n^{\mathrm{wor}}(\varepsilon,d;\mathrm{App}_d) \leq C\,\varepsilon^{-p}\,d^{q_2} \quad \text{ for all } d \in \mathbb{N} \text{ and } \varepsilon \in (0,1].$$

For any given $d \in \mathbb{N}$ we can take, say, $\varepsilon = \varepsilon(d) = \frac{1}{2} \min\{1, C_{1,d}^{-1}\}$ to conclude

$$2^{\wedge} \left(\frac{1}{2 + C_{\gamma}} \sum_{j=1}^{d} \gamma_{d,j} \right) < C' \max\{1, C_{1,d}^{p}\} d^{q_2}$$
 (4.24)

for some C'>0 independent of d. If we now assume that $C_{1,d}\in\mathcal{O}(d^{q_1})$ then the right-hand side of the last inequality belongs to $\mathcal{O}(d^{pq_1+q_2})$ as $d\to\infty$. Provided that $\max\{q_1,q_2\}>0$ this is equivalent to the boundedness of $\sum_{j=1}^d \gamma_{d,j}/\ln(d+1)$ so that we arrive at $q(\gamma)\leq 1$, as claimed.

Finally, the case of strong polynomial tractability can be treated similarly by setting $q_1=q_2=0$ in the above bounds. Then we deduce that $\sum_{j=1}^d \gamma_{d,j}$ is uniformly bounded in d, which implies $p(\gamma) \leq 1$.

Of course, the conditions $q(\gamma) \leq 1$ and $p(\gamma) \leq 1$ are also necessary for polynomial and strong polynomial tractability with respect to smaller classes of algorithms such as, e.g., $\mathcal{A}_d^{n,\text{lin}}(\Lambda^{\text{all}})$.

Observe that one of the improvements compared to [Wei12b, Theorem 2] is the possibility of choosing the uniform upper bound for the generator weights, C_{γ} , different from 1. Moreover, now we have weaker conditions on the embedding constant $C_{1,d}$. For the application we have in mind we will see that there is still $C_{1,d} = 1$. But we note in passing that the conclusions stated for (strong) polynomial tractability are only special instances of the more general bound (4.24), which we will not investigate further.

Next we assume (A4.2) and show that slightly stronger conditions on the product weights γ than in Proposition 4.6 are sufficient for polynomial and strong polynomial tractability, respectively. This is stated in the next assertion which can be found as Theorem 3 in [Wei12b].

PROPOSITION 4.7 (Sufficient conditions). Suppose that assumption (A4.2) holds true for some $t \in (0,1]$. Consider L_{∞} -approximation over $(\mathcal{F}_{d}^{\gamma})_{d \in \mathbb{N}}$ in the worst case setting with respect to the class of linear algorithms $\mathcal{A}_{d}^{n,\text{lin}}(\Lambda^{\text{all}})$ and the absolute error criterion. Then

- $q(\gamma) < t$ implies polynomial tractability,
- $p(\gamma) < t$ implies strong polynomial tractability.

Proof. Due to (A4.2), the restriction of the algorithm $A_{n,d}^*$ in Corollary 4.4 from \mathcal{H}_d^{γ} to \mathcal{F}_d^{γ} is admissible for L_{∞} -approximation over \mathcal{F}_d^{γ} . Furthermore, due to the linearity of $A_{n,d}^*$ we have

$$||f - A_{n,d}^* f \mid \mathcal{L}_{\infty}([0,1]^d)|| \leq \Delta^{\text{wor}}(A_{n,d}^*; \mathrm{id}_d \colon \mathcal{B}(\mathcal{H}_d^{\gamma}) \to \mathcal{L}_{\infty}([0,1]^d))||f \mid \mathcal{H}_d^{\gamma}||$$

$$\leq \Delta^{\text{wor}}(A_{n,d}^*; \mathrm{id}_d \colon \mathcal{B}(\mathcal{H}_d^{\gamma}) \to \mathcal{L}_{\infty}([0,1]^d))C_{2,d}||f \mid \mathcal{F}_d^{\gamma}||$$

for all $f \in \mathcal{F}_d^{\gamma}$. Therefore we can estimate the *n*th minimal error by

$$\begin{split} e^{\text{wor}}(n, d; \operatorname{App}_{d}) &\leq \Delta^{\text{wor}} \left(A_{n,d}^{*} \big|_{\mathcal{F}_{d}^{\gamma}}; \operatorname{App}_{d} \colon \mathcal{B}(\mathcal{F}_{d}^{\gamma}) \to \operatorname{L}_{\infty}([0, 1]^{d}) \right) \\ &\leq C_{2,d} \Delta^{\text{wor}}(A_{n,d}^{*}; \operatorname{id}_{d} \colon \mathcal{B}(\mathcal{H}_{d}^{\gamma}) \to \operatorname{L}_{\infty}([0, 1]^{d})) \\ &\leq a a_{\tau} \exp \left(b \sum_{j=1}^{d} \left(\gamma_{d,j} \right)^{t} + b_{\tau} \sum_{j=1}^{d} \left(\gamma_{d,j} \right)^{\tau} \right) n^{-(1-\tau)/(2\tau)}, \end{split}$$

where τ is an arbitrary number from (1/2, 1). Choosing n such that the right-hand side is not greater than a given $\varepsilon \in (0, 1]$, we obtain an estimate for the information complexity with respect to the class of linear algorithms,

$$n^{\text{wor}}(\varepsilon, d; \operatorname{App}_d) \le c_1 \varepsilon^{-2\tau/(1-\tau)} \exp\left(c_2 \sum_{j=1}^d (\gamma_{d,j})^t + c_3 \sum_{j=1}^d (\gamma_{d,j})^\tau\right), \tag{4.25}$$

where the non-negative constants c_1, c_2, c_3 only depend on τ , a and b.

Suppose that $q(\gamma) < t$. Then $Q_{\kappa}(\gamma)$ is finite for every $\kappa > q(\gamma)$. Taking $\kappa = t$ we obtain

$$\frac{\sum_{j=1}^{d} (\gamma_{d,j})^{t}}{\ln(d+1)} \ln(d+1) \le (Q_{t}(\gamma) + \delta) \ln(d+1) = \ln(d+1)^{Q_{t}(\gamma) + \delta}$$

for every $\delta > 0$ whenever d is larger than a certain $d_{\delta} \in \mathbb{N}$. This means that the factor $\exp(c_2 \sum_{j=1}^{d} (\gamma_{d,j})^t)$ in (4.25) is polynomially dependent on d. On the other hand, we can choose $\tau \in (\max\{q(\gamma), 1/2\}, 1)$ such that $Q_{\tau}(\gamma)$ is finite and thus the factor

 $\exp(c_3\sum_{j=1}^d(\gamma_{d,j})^{\tau})$ in (4.25) is polynomially dependent on d as well. So, for this value of τ we can rewrite (4.25) as

$$n^{\text{wor}}(\varepsilon, d; \text{App}_d) \in \mathcal{O}(\varepsilon^{-2\tau/(1-\tau)}(d+1)^{c_4}),$$

with c_4 , as well as the implied factor in the \mathcal{O} -notation, independent of d and ε , which means that the problem is polynomially tractable, as claimed.

Finally suppose that $p(\gamma) < t$. Then the sums $\sum_{j=1}^{d} (\gamma_{d,j})^t$ and $\sum_{j=1}^{d} (\gamma_{d,j})^{\tau}$ for $\tau \in$ $(\max\{p(\gamma), 1/2\}, 1)$ are both uniformly bounded in d. Consequently, (4.25) yields strong polynomial tractability.

The conditions in Proposition 4.7 are obviously also sufficient if we consider larger classes of algorithms such as, e.g., $\mathcal{A}_d^{n,\mathrm{cont}} \cup \mathcal{A}_d^{n,\mathrm{adapt}}$. Moreover note that the proof given above also provides explicit upper bounds for the exponents of tractability.

In the following remark we briefly discuss the different roles of assumptions (A4.1) and (A4.2).

REMARK 4.8. Assumption (A4.1) is used to find a lower bound on the information complexity for the space \mathcal{F}_d^{γ} as long as the space \mathcal{P}_d^{γ} is continuously embedded in \mathcal{F}_d^{γ} with an embedding constant which grows at most polynomially with the dimension d. Such an embedding can be shown for several different classes of functions.

In contrast, assumption (A4.2) is used to find an upper bound on the information complexity for the space \mathcal{F}_d^{γ} as long as it is continuously embedded in the unanchored weighted Sobolev space $\mathcal{H}_d^{\gamma} = \mathcal{H}(K_d^{\gamma})$ with an embedding constant depending exponentially on the sum of some power of the generators $\gamma_{d,j}$ of the product weights γ . This considerably restricts the choice of \mathcal{F}_d^{γ} . We need this assumption in order to use the linear algorithm $A_{n,d}^*$ defined on the space \mathcal{H}_d^{γ} and the error bound given in Corollary 4.4. Obviously, we can replace \mathcal{H}_d^{γ} in (A4.2) by any other space which contains at least \mathcal{P}_d^{γ} and for which we know a linear algorithm using n linear functionals whose worst case error is polynomial in n^{-1} with an explicit dependence on the product weights γ . \Box

Now we show that assumptions (A4.1) and (A4.2) allow us to characterize weak tractability and the curse of dimensionality.

THEOREM 4.9 (Weak tractability and the curse of dimensionality). Suppose that for a sequence $(\mathcal{F}_d^{\gamma})_{d\in\mathbb{N}}$ of Banach spaces equipped with product weights γ assumptions (A4.1) and (A4.2) hold true with some parameter $t \in (0,1]$. Consider the L_{∞} -approximation problem App in the worst case setting and with respect to the absolute error criterion. Then the following statements are equivalent:

- (i) The problem is weakly tractable with respect to the class A_d^{n,lin}(Λ^{all}).
 (ii) The problem is weakly tractable with respect to the class A_d^{n,cont} ∪ A_d^{n,adapt}.
- (iii) There is no curse of dimensionality for the class $\mathcal{A}_d^{n,\text{lin}}(\Lambda^{\text{all}})$.
- (iv) There is no curse of dimensionality for the class $\mathcal{A}_d^{n,\mathrm{cont}} \cup \mathcal{A}_d^{n,\mathrm{adapt}}$.
- (v) For all $\kappa > 0$ we have $\lim_{d\to\infty} \frac{1}{d} \sum_{j=1}^d (\gamma_{d,j})^{\kappa} = 0$.
- (vi) There exists $\kappa \in (0,t)$ such that $\lim_{d\to\infty} \frac{1}{d} \sum_{j=1}^d (\gamma_{d,j})^{\kappa} = 0$.

Proof. We start by showing that (vi) implies (i), i.e.,

$$\lim_{\varepsilon^{-1}+d\to\infty}\frac{\ln n^{\mathrm{wor}}(\varepsilon,d;\mathrm{App}_d)}{\varepsilon^{-1}+d}=0,$$

where the information complexity is taken with respect to the class $\mathcal{A}_d^{n,\text{lin}}(\Lambda^{\text{all}})$ of linear algorithms that use continuous linear functionals. By the arguments used in the proof of Proposition 4.7 we obtain estimate (4.25) for all ε in (0, 1], as well as for every $d \in \mathbb{N}$, and all $\tau \in (1/2, 1)$, due to assumption (A4.2). Clearly, for $\kappa \in (0, t)$ as in the hypothesis and $t \in (0, 1]$ as in the embedding condition, we find $\tau \in (1/2, 1)$ such that $\kappa < \min\{t, \tau\}$. So, since $\gamma_{d,j} \leq C_{\gamma}$, we can estimate

$$\sum_{j=1}^{d} (\gamma_{d,j})^s = C_{\gamma}^s \sum_{j=1}^{d} \left(\frac{\gamma_{d,j}}{C_{\gamma}}\right)^s \le C_{\gamma}^{s-\kappa} \sum_{j=1}^{d} (\gamma_{d,j})^{\kappa} \le C \sum_{j=1}^{d} (\gamma_{d,j})^{\kappa},$$

where s equals either t or τ , and $C = \max\{1, C_{\gamma}\}$. Therefore the right-hand side of (4.25) can be estimated from above and thus

$$\frac{\ln n^{\text{wor}}(\varepsilon, d; \operatorname{App}_d)}{\varepsilon^{-1} + d} \le \frac{\ln c_1}{\varepsilon^{-1} + d} + \frac{2\tau}{1 - \tau} \frac{\ln \varepsilon^{-1}}{\varepsilon^{-1} + d} + C \max\{c_2, c_3\} \frac{\sum_{j=1}^d (\gamma_{d,j})^{\kappa}}{\varepsilon^{-1} + d},$$

which tends to zero when $\varepsilon^{-1} + d \to \infty$, as claimed.

Clearly, (i) \Rightarrow (ii) \Rightarrow (iv) and (i) \Rightarrow (iii) \Rightarrow (iv). Moreover, the implication from (v) to (vi) is obvious. Hence, it only remains to show that (iv) \Rightarrow (v).

From (A4.1) we have estimate (4.23). Then the absence of the curse of dimensionality implies

$$\lim_{d \to \infty} \frac{1}{d} \sum_{j=1}^{d} \gamma_{d,j} = 0.$$

Now Jensen's inequality yields

$$\frac{1}{d} \sum_{i=1}^{d} \gamma_{d,j} \ge \left(\frac{1}{d} \sum_{i=1}^{d} (\gamma_{d,j})^{\kappa}\right)^{1/\kappa} \quad \text{for } 0 < \kappa \le 1,$$

because $f(y) = y^{\kappa}$ is a concave function for y > 0. This shows

$$\lim_{d \to \infty} \frac{1}{d} \sum_{j=1}^{d} (\gamma_{d,j})^{\kappa} = 0 \quad \text{for all } 0 < \kappa \le 1.$$

Finally, for every $\kappa \geq 1$ we can estimate $\gamma_{d,j} \geq C_{\gamma}^{1-\kappa} (\gamma_{d,j})^{\kappa}$ since $\gamma_{d,j} \leq C_{\gamma}$ for $j = 1, \ldots, d$. Therefore $\lim_{d \to \infty} d^{-1} \sum_{j=1}^{d} (\gamma_{d,j})^{\kappa} = 0$ also holds true for $\kappa > 1$, and the proof of Theorem 4.9 is complete.

4.3.3. Conclusions and applications. In this last part of Section 4.3 we give some examples to illustrate the complexity results obtained. To this end, we only have to prove the corresponding embeddings, i.e. we need to verify assumption (A4.1) and/or (A4.2) from the beginning of Section 4.3.2.

EXAMPLE 4.10 (Limiting cases \mathcal{P}_d^{γ} and \mathcal{H}_d^{γ}). To begin with, we check the case where $\mathcal{F}_d^{\gamma} = \mathcal{P}_d^{\gamma}$ for every $d \in \mathbb{N}$. Then (A4.1) obviously holds with $C_{1,d} = 1$, i.e. c = 1 and

 $q_1=0$. To prove (A4.2), note that the algebraical inclusion $\mathcal{F}_d^{\gamma}\subset\mathcal{H}_d^{\gamma}$ is trivial by the definition of $\mathcal{H}_d^{\gamma}=\mathcal{H}(K_d^{\gamma})$ given in Section 4.1.2. For $f\in\mathcal{F}_d^{\gamma}=\mathcal{P}_d^{\gamma}$ we calculate

$$||f| \mathcal{H}_{d}^{\gamma}||^{2} \leq \sum_{\alpha \in \{0,1\}^{d}} \frac{1}{\gamma_{\alpha}} \int_{[0,1]^{d}} ||D^{\alpha}f| L_{\infty}([0,1]^{d})||^{2} d\lambda^{d}(\boldsymbol{x}) \leq ||f| \mathcal{F}_{d}^{\gamma}||^{2} \sum_{\alpha \in \{0,1\}^{d}} \gamma_{\alpha}$$

using (4.10), as well as (4.17). Hence the norm of the embedding $\mathcal{F}_d^{\gamma} \hookrightarrow \mathcal{H}_d^{\gamma}$ is bounded by

$$\left(\sum_{\alpha \in \{0,1\}^d} \gamma_{\alpha}\right)^{1/2} = \left(\prod_{j=1}^d (1 + \gamma_{d,j})\right)^{1/2} \le \exp\left(\frac{1}{2} \sum_{j=1}^d \gamma_{d,j}\right).$$

So, with $a=1,\ b=1/2$, and t=1, assumption (A4.2) is fulfilled as well, and we can apply the assertions stated in Section 4.3.2 for the spaces $\mathcal{F}_d^{\gamma} = \mathcal{P}_d^{\gamma},\ d \in \mathbb{N}$.

We now turn to the case $\mathcal{F}_d^{\gamma} = \mathcal{H}_d^{\gamma}$. Unfortunately, the above estimate indicates that (A4.1) may not hold for $\mathcal{F}_d^{\gamma} = \mathcal{H}_d^{\gamma}$ with $C_{1,d} \in \mathcal{O}(d^{q_1})$ without imposing additional conditions on the product weights γ . Nevertheless, in this case assumption (A4.2) is trivially true with $C_{2,d} = 1$, i.e., a = 1, b = 0, and t = 1. Therefore we can apply Proposition 4.7 for this space. Thus the problem is polynomially tractable if $q(\gamma) < 1$, and we have strong polynomial tractability if $p(\gamma) < 1$. It can be shown that these conditions are also necessary (see Section 4.4). \square

Next we discuss a more advanced sequence of Banach function spaces.

Example 4.11 $(C^{(1,\ldots,1)})$. For every $d \in \mathbb{N}$ consider the space

$$\mathcal{F}_d^{\gamma} = \{ f : [0,1]^d \to \mathbb{R} \mid f \in C^{(1,\dots,1)}([0,1]^d), \text{ where } ||f| \mid \mathcal{F}_d^{\gamma}|| < \infty \}$$

of functions which are once continuously differentiable in every coordinate direction, where

$$||f| \mathcal{F}_d^{\gamma}|| = \max_{\alpha \in \{0,1\}^d} \frac{1}{\gamma_{\alpha}} ||D^{\alpha} f| L_{\infty}([0,1]^d)||.$$

Since \mathcal{P}_d^{γ} is a linear subset of \mathcal{F}_d^{γ} and, due to (4.17), the norm $\|\cdot \mid \mathcal{P}_d^{\gamma}\|$ is simply the restriction of $\|\cdot \mid \mathcal{F}_d^{\gamma}\|$ we have $\mathcal{P}_d^{\gamma} \hookrightarrow \mathcal{F}_d^{\gamma}$ with an embedding factor $C_{1,d} = 1$ and hence (A4.1) holds true. For the norm $C_{2,d}$ of the embedding $\mathcal{F}_d^{\gamma} \hookrightarrow \mathcal{H}_d^{\gamma}$, the same estimates hold exactly as in the previous example and, moreover, the set inclusion is obvious. Therefore also assumption (A4.2) is fulfilled and we can apply the propositions and theorem of Section 4.3.2 to the sequence $(\mathcal{F}_d^{\gamma})_{d\in\mathbb{N}}$. \square

Our last example $\mathcal{F}_d^{\gamma} = F_d^{\gamma}$, for all $d \in \mathbb{N}$, finally shows that even very high smoothness does not improve the conditions for tractability.

Example 4.12 (C^{∞}) . For $d \in \mathbb{N}$ and product weights γ let

$$\mathcal{F}_{d}^{\gamma} = F_{d}^{\gamma} = \{ f : [0, 1]^{d} \to \mathbb{R} \mid f \in C^{\infty}([0, 1]^{d}) \text{ with } ||f| ||F_{d}^{\gamma}|| < \infty \},$$

where the norm is given by (4.5). Obviously, $\mathcal{P}_d^{\gamma} \subset C^{\infty}$, because functions from \mathcal{P}_d^{γ} are at most linear in each coordinate. This moreover implies that $D^{\alpha}f \equiv 0$ for all α in $\mathbb{N}_0^d \setminus \{0,1\}^d$. Therefore, once again we have

$$||f| \mathcal{P}_d^{\gamma}|| = \max_{\alpha \in \{0,1\}^d} \frac{1}{\gamma_{\alpha}} ||D^{\alpha}f| L_{\infty}([0,1]^d)|| = ||f| \mathcal{F}_d^{\gamma}|| \quad \text{for all } f \in \mathcal{P}_d^{\gamma}.$$

Together this yields $\mathcal{P}_d^{\gamma} \hookrightarrow \mathcal{F}_d^{\gamma}$ with an embedding constant $C_{1,d} = 1$ for all $d \in \mathbb{N}$. In addition, also (A4.2) can be verified as in the examples above. So, even infinite smoothness leads to the same conditions for tractability and the curse of dimensionality as before. \square

Note that in Example 4.12 we do not need to claim a product structure for the weights according to multi-indices $\alpha \in \mathbb{N}_0^d \setminus \{0,1\}^d$. Furthermore, the space F_d^{γ} is a generalization of the space F_d studied in [NW09]. For $\gamma_{\alpha} \equiv 1$ we reproduce the intractability result stated there because then F_d^{γ} equals F_d for each $d \in \mathbb{N}$.

In conclusion we discuss the tractability behavior of uniform approximation defined on one of the spaces \mathcal{F}_d^{γ} above using a special class of product weights γ which are independent of the dimension d. That is, for the generator weights we claim that

$$\gamma_{d,j} \equiv \gamma^{(j)} \in \Theta(j^{-\beta}) \quad \text{for some } \beta \ge 0$$
 (4.26)

and all j and $d \in \mathbb{N}$. The polynomial behavior of $\gamma^{(j)}$ imposed is a typical example in the theory of product weights. Clearly, $p(\gamma)$ is finite if and only if $\beta > 0$, and if so then $p(\gamma) = 1/\beta$. For details see [NW08, Section 5.3.4].

If $\beta = 0$ then the L_∞-approximation problem App = $(\mathrm{App}_d)_{d \in \mathbb{N}}$ is intractable (more precisely it suffers from the curse of dimensionality) due to Theorem 4.9(v), since then $d^{-1} \sum_{j=1}^{d} \gamma_{d,j}$ does not tend to 0. For $\beta \in (0,1)$ an easy computation yields $q(\gamma) > 1$. So, using Proposition 4.6 we conclude polynomial intractability in this case. On the other hand, we have for all δ and κ with $0 < \delta < \kappa \le 1$,

$$\frac{\sum_{j=1}^{d} j^{-\kappa}}{d} = \frac{\sum_{j=1}^{d} j^{-\kappa} d^{\kappa - (1+\delta)}}{d^{\kappa - \delta}} \le \frac{\sum_{j=1}^{d} j^{-(1+\delta)}}{d^{\kappa - \delta}} \to 0 \quad \text{for } d \to \infty$$

and if $\kappa > 1$ then the leftmost fraction obviously tends to zero, too. Hence condition (vi) of Theorem 4.9 holds and the problem is weakly tractable for all $\beta > 0$.

For $\beta = 1$ we use inequality (4.23) from Proposition 4.6 and estimate

$$\sum_{i=1}^{d} \gamma_{d,j} \ge c \ln(d+1)$$

for some positive c. Therefore, for all $\varepsilon \in (0,1)$ the information complexity $n(\varepsilon,d;\operatorname{App}_d)$ is lower bounded polynomially in $d \in \mathbb{N}$. This proves that strong polynomial tractability does not hold for $\beta=1$. Moreover, it is easy to show that in this case the sufficient condition $q(\gamma) < 1$ for polynomial tractability is not fulfilled. So, we do not know whether polynomial tractability holds or not.

Finally, consider $\beta > 1$ in (4.26). Then we easily see that $p(\gamma) = 1/\beta < 1 = t$. Thus Proposition 4.7 provides strong polynomial tractability in this situation.

In summary, we proved all the assertions claimed at the end of Section 4.1.1.

4.4. Possible extensions and further results. Note that the main result of this chapter, the lower bound given in Theorem 4.5, can be easily transferred from $[0,1]^d$ to more general domains $\Omega_d \subset \mathbb{R}^d$. Indeed, the case $\Omega_d = [c_1, c_2]^d$, where $c_1 < c_2$, can be immediately obtained using the techniques presented above. It turns out that in this case

we have to modify estimate (4.18) by a constant which depends only on the length of the interval $[c_1, c_2]$. Consequently, the general tractability behavior does not change.

Another extension of the results obtained is possible if we consider L_p -norms (with $1 \le p < \infty$) instead of the L_{∞} -norm. In Section 4.4.1 we briefly discuss these norms for the unweighted case. Then the modifications for the weighted case are obvious and thus we leave it for the interested reader. In passing we correct a small mistake of [NW09].

Finally, in Section 4.4.2, we show that the algorithm studied in Corollary 4.4 is essentially optimal for the uniform approximation problem on the unanchored weighted Sobolev space $\mathcal{H}(K_d^{\gamma})$ defined in Example 4.1.

4.4.1. L_p -approximation. As in [Wei12b, Section 7] we follow Novak and Woźniakowski [NW09] and define the spaces

$$F_{d,p} = \left\{ f \in C^{\infty}([c_1, c_2]^d) \mid ||f| \mid F_{d,p}|| = \sup_{\alpha \in \mathbb{N}_a^d} ||D^{\alpha} f| \mid L_p([c_1, c_2]^d)|| < \infty \right\}$$

for $1 \leq p < \infty$ and $d \in \mathbb{N}$, where we assume that $l = c_2 - c_1 > 0$. In what follows we want to approximate $f \in F_{d,p}$ in the norm of L_p . That is, we modify (4.2) and consider the problem $S^p = (S_d^p)_{d \in \mathbb{N}}$ given by

$$S_d^p = \mathrm{id}_d^p \colon \mathcal{B}(F_{d,p}) \to \mathrm{L}_p([c_1, c_2]^d), \quad f \mapsto \mathrm{id}_d^p(f) = f.$$

Hence we analyze the nth minimal worst case error

$$e_p^{\text{wor}}(n, d; \text{id}_d^p) = \inf_{A_{n,d}} \sup_{f \in \mathcal{B}(F_{d,p})} ||f - A_{n,d}(f)|| L_p([c_1, c_2]^d)||,$$

which now depends on the additional integrability parameter p. Observe that, without loss of generality, we can restrict ourselves to the case $[c_1, c_2] = [0, l]$.

In order to derive a lower bound analogue to (4.3) and Theorem 4.5, i.e.,

$$e_n^{\text{wor}}(n, d; \text{id}_d^p) \ge 1$$
 for $n < 2^s$,

we once again use Proposition 2.2 with $\mathcal{F} = F_{d,p}$ and $\mathcal{G} = \mathcal{L}_p([0,l]^d)$ (²⁶). The authors of [NW09] suggest to use the subspace $V_d^{(k)} \subset F_{d,p}$ defined as

$$V_d^{(k)} = \operatorname{span} \Big\{ g_i \colon [0, l]^d \to \mathbb{R}, \, \boldsymbol{x} \mapsto g_{\boldsymbol{i}}(\boldsymbol{x}) = \prod_{j=1}^s \Big(\sum_{m=(j-1)k+1}^{jk} x_m \Big)^{i_j} \, \, \Big| \, \, \boldsymbol{i} \in \{0, 1\}^s \Big\},$$

where $s = \lfloor d/k \rfloor$ and $k \in \mathbb{N}$ is such that $kl \geq 2(p+1)^{1/p}$. Hence, if $l < 2(p+1)^{1/p}$ then we have to use blocks of variables with size k > 1 in order to guarantee (2.1), that is, to fulfill the condition

$$||g| F_{d,p}|| \le ||g| L_p([0,l]^d)||$$
 for all $g \in V_d^{(k)}$. (4.27)

Therefore Novak and Woźniakowski defined $k = \lceil 2(p+1)^{1/p}/l \rceil$, but this is too small as the following example shows.

EXAMPLE 4.13. For $d \ge 4$ take l = 1, i.e. $[c_1, c_2]^d = [0, 1]^d$, and p = 1. Then k = 4 should be a proper choice, but for $g^*(\mathbf{x}) = (x_1 + x_2 + x_3 + x_4) - 2$ it can be checked (using a

 $[\]binom{26}{}$ Note that it is sufficient to restrict ourselves to the case r=1 since now we do not need to take care of embedding constants as in the proof of Proposition 4.6.

computer algebra system) that

$$||g^*| L_1([0,1]^d)|| = \frac{7}{15} < 1 = \left\| \frac{\partial g^*}{\partial x_1} | L_1([0,1]^d) \right\|.$$

This obviously contradicts (4.27). \Box

For an exhaustive proof that a slightly larger choice of $k \in \mathbb{N}$ suffices to obtain the desired intractability result we need to show the following technical lemma first. Its proof is based on some well-known arguments from Banach space geometry.

LEMMA 4.14. Let $p \in [1, \infty)$ and $k \in \mathbb{N}$. Then

$$\mathfrak{I}_{k,p} = \int_{[-1/2,1/2]^k} \left| \sum_{m=1}^k z_m \right|^p d\lambda^k(z) \ge C_p k^{p/2}$$
(4.28)

with some $C_p \geq 1/[(2\sqrt{2})^p(1+p)]$ independent of k.

Proof. For k=1 we easily calculate $\mathfrak{I}_{1,p}=1/[2^p(1+p)]$. Hence, without loss of generality we can assume $k\geq 2$ in what follows.

To abbreviate the notation, let us define

$$f = f_k \colon \mathbb{R}^k \to \mathbb{R}, \quad \boldsymbol{z} = (z_1, \dots, z_k) \mapsto f(\boldsymbol{z}) = \sum_{m=1}^k z_m,$$
 (4.29)

for any fixed $k \geq 2$. Moreover, for given vectors $\boldsymbol{z}, \boldsymbol{y} \in \mathbb{R}^k$, let $\langle \boldsymbol{z}, \boldsymbol{y} \rangle$ denote the inner product $\sum_{m=1}^k z_m y_m$ in \mathbb{R}^k . In the special case $\boldsymbol{y} = \boldsymbol{\xi} = (1/\sqrt{k})(1, \dots, 1) \in \mathbb{S}^{k-1}$ we have $\langle \boldsymbol{z}, \boldsymbol{\xi} \rangle = t$ for a given $t \in \mathbb{R}$ if and only if $f(\boldsymbol{z}) = t\sqrt{k}$. Furthermore note that every \boldsymbol{y} in the k-dimensional unit sphere $\mathbb{S}^{k-1} \subset \mathbb{R}^k$ uniquely defines a hyperplane $\boldsymbol{y}^{\perp} = \{\boldsymbol{z} \in \mathbb{R}^k \mid \langle \boldsymbol{z}, \boldsymbol{y} \rangle = 0\}$ perpendicular to \boldsymbol{y} which contains zero. Therefore, for $\boldsymbol{y} = \boldsymbol{\xi}$ and every $t \in [0, \infty)$, the set

$$\mathfrak{H}_t = \boldsymbol{\xi}^{\perp} + t \boldsymbol{\xi} = \{ \boldsymbol{z} \in \mathbb{R}^k \mid \langle \boldsymbol{z}, \boldsymbol{\xi} \rangle = t \}$$

describes a parallel shifted hyperplane in \mathbb{R}^k with distance t to the origin. Using Fubini's theorem, this leads to the following representation:

$$\begin{split} \mathfrak{I}_{k,p} &= \int_{[-1/2,1/2]^k} \left| f(\boldsymbol{z}) \right|^p \, \mathrm{d}\lambda^k(\boldsymbol{z}) = 2 \int_{[-1/2,1/2]^k} f(\boldsymbol{z})^p \, \mathrm{d}\lambda^k(\boldsymbol{z}) \\ &= 2k^{p/2} \int_0^\infty t^p \left(\int_{[-1/2,1/2]^k \cap \mathfrak{H}_t} 1 \, \mathrm{d}\lambda^k(\boldsymbol{z}) \right) \mathrm{d}\lambda^1(t). \end{split}$$

Now we see that the inner integral gives the (k-1)-dimensional volume

$$v(t) = \lambda^{k-1}([-1/2, 1/2]^k \cap \mathfrak{H}_t)$$

of the parallel section of the unit cube with the hyperplanes defined above. By Ball's famous theorem we know that $v(0) \leq \sqrt{2}$ independently of k; see, e.g., Chapter 7 in the monograph of Koldobsky [Kol05]. Moreover taking $\mathfrak{H}_0 = \boldsymbol{\xi}^{\perp}$ provides a central hyperplane section of the unit cube. From this observation we conclude that

$$\int_0^\infty v(t) \, \mathrm{d}\lambda^1(t) = \frac{1}{2} \lambda^k([-1/2, 1/2]^k) = \frac{1}{2}$$

by the symmetry of $[-1/2, 1/2]^k$ with respect to \mathfrak{H}_0 . In addition, by Brunn's theorem (cf. [Kol05, Theorem 2.3]), the function v is non-negative and non-increasing on $[0, \infty)$. Thus v is related to the distribution function of a certain non-negative real-valued random variable X up to some normalizing factor, i.e. $v(t) = v(0)\mathbb{P}(\{X \geq t\})$. Using Hölder's inequality (27) we obtain $\mathbb{E}(X^{1+p}) \geq (\mathbb{E}X)^{1+p}$ and, respectively,

$$\Im_{k,p} = k^{p/2} \cdot 2 \int_0^\infty t^p \, v(t) \, \mathrm{d}\lambda^1(t) \ge k^{p/2} \frac{2}{v(0)^p \, (1+p)} \left(\int_0^\infty v(t) \, \mathrm{d}\lambda^1(t) \right)^{1+p}$$

by integration by parts.

In summary we have shown (4.28) and hence the proof is complete. \blacksquare

Now the intractability result mentioned at the beginning of this section reads as follows:

Proposition 4.15. Let $1 \le p < \infty$ and l > 0. Moreover, choose $k \in \mathbb{N}$ such that

$$k \ge \kappa_{p,l} = \lceil 8(p+1)^{2/p}/l^2 \rceil.$$
 (4.30)

Then condition (4.27) holds for $V_d^{(k)} \subset F_{d,p}$. Therefore the L_p -approximation problem $S^p = (\operatorname{id}_d^p : F_{d,p} \to L_p([0,l]^d))_{d \in \mathbb{N}}$ suffers from the curse of dimensionality since

$$e_n^{\text{wor}}(n, d; \mathrm{id}_d^p) \ge 1$$
 for all $n < 2^{\lfloor d/k \rfloor}$

and every $d \in \mathbb{N}$.

Proof. Due to the structure of the functions g from $V_d^{(k)}$, it suffices to show that

$$||D^{\boldsymbol{\alpha}}g| L_p([0,l]^{ks})|| \le ||g| L_p([0,l]^{ks})||$$
 for all $g \in V_d^{(k)}$ and every $\boldsymbol{\alpha} \in \mathbb{M}_d^{(k)}$,

where the set of multi-indices $\mathbb{M}_d^{(k)}$ is defined by

$$\mathbb{M}_d^{(k)} = \left\{ \boldsymbol{\alpha} = (\alpha_1, \dots, \alpha_{ks}) \in \{0, 1\}^{ks} \mid \sum_{m \in I_j} \alpha_m \le 1 \text{ for all } j = 1, \dots, s \right\}$$

and $I_j = \{(j-1)k+1, \ldots, jk\}$. Observe that $\mathbb{M}_d^{(k)}$ depends on d via $s = \lfloor d/k \rfloor$. Similar to the proof of Theorem 4.5, we only need to consider the case $\boldsymbol{\alpha} = \boldsymbol{e_t} \in \{0,1\}^{ks}$ with $t \in I_j$. The rest then follows by induction.

Given $t \in I_j$ for some $j \in \{1, ..., s\}$ we can represent every fixed $g \in V_d^{(k)}$, as well as its partial derivative $D^{e_t}g$, by some functions $a, b : [0, l]^{k(s-1)} \to \mathbb{R}$ (depending on g and j) such that

$$g(\boldsymbol{x}) = a(\widetilde{\boldsymbol{x}}) \sum_{m=1}^{k} y_m + b(\widetilde{\boldsymbol{x}})$$
 and $(D^{\boldsymbol{e_t}}g)(\boldsymbol{x}) = a(\widetilde{\boldsymbol{x}}), \quad \boldsymbol{x} \in [0, l]^{ks}.$

Here we split the ks-dimensional vector $\boldsymbol{x}=(\boldsymbol{x}_{I_1},\ldots,\boldsymbol{x}_{I_{j-1}},\boldsymbol{y},\boldsymbol{x}_{I_{j+1}},\ldots,\boldsymbol{x}_{I_s})$ into $\widetilde{\boldsymbol{x}}=(\boldsymbol{x}_{I_1},\ldots,\boldsymbol{x}_{I_{j-1}},\boldsymbol{x}_{I_{j+1}},\ldots,\boldsymbol{x}_{I_s})\in[0,l]^{k(s-1)}$ and $\boldsymbol{y}=(y_1,\ldots,y_k)\in[0,l]^k$, where \boldsymbol{x}_{I_j} denotes the k-dimensional block of components x_m in \boldsymbol{x} with coordinates $m\in I_j$. Using this

 $^(^{27})$ See also [Kol05, Lemma 7.5].

representation we can rewrite the inequality $||D^{e_t}g| L_p([0,l]^{ks})|| \le ||g| L_p([0,l]^{ks})||$ as

$$\int_{[0,l]^{k(s-1)}} \int_{[0,l]^k} |a(\widetilde{\boldsymbol{x}})|^p d\lambda^k(\boldsymbol{y}) d\lambda^{k(s-1)}(\widetilde{\boldsymbol{x}})
\leq \int_{[0,l]^{k(s-1)}} \int_{[0,l]^k} |a(\widetilde{\boldsymbol{x}}) \sum_{m=1}^k y_m + b(\widetilde{\boldsymbol{x}})|^p d\lambda^k(\boldsymbol{y}) d\lambda^{k(s-1)}(\widetilde{\boldsymbol{x}})$$

so that it is enough to prove a pointwise estimate of the inner integrals for $(\lambda^{k(s-1)}$ -almost every) fixed $\tilde{x} \in [0, l]^{k(s-1)}$ with $a = a(\tilde{x}) \neq 0$. A simple calculation yields

$$\int_{[0,l]^k} \left| a \sum_{m=1}^k y_m + b \right|^p d\lambda^k(\boldsymbol{y}) = l^{p+k} \int_{[-1/2,1/2]^k} \left| a \sum_{m=1}^k z_m + b' \right|^p d\lambda^k(\boldsymbol{z})$$

for some constant $b' \in \mathbb{R}$ that depends on $b = b(\widetilde{x})$. Note that the right-hand side of the last equality is minimized for b' = 0. Therefore we can estimate the left-hand side from below by

$$\int_{[0,l]^k} \left| a \sum_{m=1}^k y_m + b \right|^p d\lambda^k(\boldsymbol{y}) \ge |a|^p l^{p+k} \int_{[-1/2,1/2]^k} \left| \sum_{m=1}^k z_m \right|^p d\lambda^k(\boldsymbol{z})
= \int_{[0,l]^k} |a|^p d\lambda^k(\boldsymbol{y}) l^p \int_{[-1/2,1/2]^k} \left| \sum_{m=1}^k z_m \right|^p d\lambda^k(\boldsymbol{z}).$$

To complete the proof we are left with showing that our choice of $k \geq \kappa_{p,l}$, with $\kappa_{p,l}$ given in (4.30), implies that

$$\int_{[-1/2,1/2]^k} \left| \sum_{m=1}^k z_m \right|^p d\lambda^k(\mathbf{z}) \ge l^{-p}; \tag{4.31}$$

but this easily follows from Lemma 4.14 above.

Actually, using other proof methods we can slightly improve the lower bound for C_p in Lemma 4.14 and thus also $\kappa_{p,l}$ in formula (4.30) of Proposition 4.15. This is the subject of our final remark in this subsection:

REMARK 4.16. Let $\mathbf{Y} = (Y_1, \dots, Y_k)$ denote a random vector of $k \in \mathbb{N}$ independent copies of some uniformly [-1/2, 1/2]-distributed random variable Y_0 . Then $\mathfrak{I}_{k,p}$ can be interpreted as the pth absolute moment $\mathbb{E}(|f_k(\mathbf{Y})|^p)$ of $f_k(\mathbf{Y})$, where f_k is again given by (4.29). In the case of even p = 2N, $N \in \mathbb{N}$, this can be calculated exactly using the multinomial theorem. For $k, N \in \mathbb{N}$ we obtain

$$\mathfrak{I}_{k,2N} = \mathbb{E}(|f_k(\boldsymbol{Y})|^{2N}) = 2^{-2N} \sum_{\substack{\boldsymbol{j} = (j_1, \dots, j_k) \in \mathbb{N}_0^k \\ j_1 + \dots + j_k = N}} {2j_1, \dots, 2j_k} \prod_{m=1}^k \frac{1}{2j_m + 1},$$

where we used the independence of the Y_m 's and the fact that

$$\mathbb{E}(Y_0^n) = \int_{-1/2}^{1/2} y^n \, \mathrm{d}\lambda^1(y) = \begin{cases} 0 & \text{if } n = 2j+1, \\ (2j+1)^{-1} \cdot 2^{-2j} & \text{if } n = 2j, \end{cases}$$

with $j \in \mathbb{N}_0$. In particular, we conclude that

$$\mathfrak{I}_{k,2} = \frac{1}{2^2 \cdot 3} k$$
 and $\mathfrak{I}_{k,4} = \frac{1}{48} k \left(k - \frac{2}{5} \right) \ge \frac{1}{2^4 \cdot 5} k^2$.

Since $\mathfrak{I}_{k,p} = \|f_k \mid \mathrm{L}_p([-1/2,1/2]^k)\|^p$ we can use the monotonicity of the Lebesgue spaces in order to estimate C_p for the remaining powers p. For $k \in \mathbb{N}$ and $1 \le q \le p < \infty$ we obtain $\mathfrak{I}_{k,p} \ge (\mathfrak{I}_{k,q})^{p/q} \ge (C_q)^{p/q} \, k^{p/2}$, i.e. $C_p \ge (C_q)^{p/q}$, provided that $\mathfrak{I}_{k,q} \ge C_q \, k^{q/2}$. Consequently, we can take

$$k \ge \begin{cases} \lceil 12/l^2 \rceil & \text{if } 2 \le p < 4, \\ \lceil 4\sqrt{5}/l^2 \rceil & \text{if } 4 \le p \end{cases}$$

to fulfill (4.31) in the proof of Proposition 4.15. This clearly improves the bound $k \geq \kappa_{p,l}$ in (4.30). \Box

Nevertheless, we want to stress that also with these improvements the lower bounds on k are not sharp since we know from [NW09] that in the limit case $p=\infty$ we can take $k=\lceil 2/l \rceil$. On the other hand, Hoeffding's inequality implies the existence of some universal constants C_p' such that $\mathfrak{I}_{k,p} \leq C_p' \, k^{p/2}$ for all $p \in [1,\infty)$ and every $k \in \mathbb{N}$. Thus the estimates on the integrals $\mathfrak{I}_{k,p}$ are of the right order in k, so that we need other proof techniques to obtain a better dependence of k on $l=c_2-c_1$.

4.4.2. Uniform approximation in the weighted Sobolev space. To show that the linear algorithm $A_{n,d}^*$ studied in Corollary 4.4 is essentially optimal for L_{∞} -approximation on the unanchored Sobolev space $\mathcal{H}_d^{\gamma} = \mathcal{H}(K_d^{\gamma})$ in the worst case setting, we study (weighted) L_2 -approximation on a related Banach space \mathcal{F}_d in the average case setting (see Example 2.15 for details). The relation of these two problems is given by the assertion below which follows from [KWW08, Theorem 1].

PROPOSITION 4.17. For $d \in \mathbb{N}$ let $\mathcal{H}(K_d)$ denote a RKHS induced by a kernel K_d : $[0,1]^d \times [0,1]^d \to \mathbb{R}$ that satisfies (4.13) (28). Moreover, define the set of non-vanishing probability density functions ρ on the unit cube by

$$\mathcal{D}_d = \bigg\{ \varrho \colon [0,1]^d \to [0,\infty) \; \bigg| \; \int_{[0,1]^d} \varrho(\boldsymbol{x}) \, \mathrm{d}\lambda^d(\boldsymbol{x}) = 1 \; \; and \; \; \varrho > 0 \; \; (\lambda^d \text{-}a.e.) \bigg\}.$$

Then, for every $n \in \mathbb{N}_0$ and all $d \in \mathbb{N}$,

$$e^{\operatorname{wor}}(n,d;\operatorname{id}_d\colon \mathcal{B}(\mathcal{H}(K_d))\to \operatorname{L}_{\infty}([0,1]^d))\geq \sup_{\varrho\in\mathcal{D}_d}e^{\operatorname{avg}}(n,d;\operatorname{id}_d^{\varrho}\colon\mathcal{F}_d\to\operatorname{L}_2^{\varrho}([0,1]^d)).$$

Here the nth minimal errors are taken with respect to all algorithms from the class $\mathcal{A}_d^{n,\text{lin}}(\Lambda^{\text{all}})$.

In particular, it follows that the (nth minimal) worst case error for L_{∞} -approximation on the unit ball of the Sobolev space \mathcal{H}_d^{γ} is lower bounded by the average case error of unweighted L_2 -approximation on the corresponding Banach space. That is, we set $K_d = K_d^{\gamma}$ and $\varrho = \chi_{[0,1]^d} \in \mathcal{D}_d$ in the following.

In turn we have (strong) polynomial tractability for the uniform approximation problem with respect to the worst case setting only if the average case L₂-approximation is polynomially tractable, as long as we consider the absolute error criterion. Due to

^{(&}lt;sup>28</sup>) Note that (4.13) clearly implies that $\int_{[0,1]^d} K_d(\boldsymbol{x},\boldsymbol{x}) \,\varrho(\boldsymbol{x}) \,\mathrm{d}\lambda^d(\boldsymbol{x})$ is finite for every probability density function ϱ on $[0,1]^d$.

[NW08, Theorem 6.1] we know that the latter holds true if and only if there exist a positive constant c_1 , non-negative q_1 , q_2 and $\tau \in (0,1)$ such that

$$c_2 = \sup_{d \in \mathbb{N}} \frac{1}{d^{q_2}} \left(\sum_{i=\lceil c_1 d^{q_1} \rceil}^{\infty} (\lambda_{d,i})^{\tau} \right)^{1/\tau} < \infty,$$

where $(\lambda_{d,i})_{i=1}^{\infty}$ denotes the sequence of eigenvalues of the correlation operator C_{ν_d} with respect to a non-increasing ordering. Moreover we have strong polynomial tractability if and only if this holds with $q_1 = q_2 = 0$.

By the observation at the end of Example 2.15 it suffices to consider the eigenvalues of $W_d^{\gamma} = (S_d^{\gamma})^{\dagger} S_d^{\gamma} : \mathcal{H}_d^{\gamma} \to \mathcal{H}_d^{\gamma}$, where the operator S_d^{γ} describes the embedding $\mathcal{H}_d^{\gamma} \hookrightarrow L_2([0,1]^d)$. Recall that these eigenvalues are given by

$$\left\{\widetilde{\lambda}_{d,\gamma,\boldsymbol{m}} = \prod_{k=1}^{d} \lambda_{1,\gamma_{d,k},m_k} = \prod_{k=1}^{d} \frac{\gamma_{d,k}}{\gamma_{d,k} + \pi^2 (m_k - 1)^2} \mid \boldsymbol{m} = (m_1, \dots, m_d) \in \mathbb{N}^d\right\}$$

(see (4.12) at the end of Example 4.1). Thus we only need to reorder this set appropriately using a rearrangement $\psi_d \colon \mathbb{N} \to \mathbb{N}^d$ such that

$$\lambda_{d,i} = \widetilde{\lambda}_{d,\gamma,\psi_d(i)} \ge \widetilde{\lambda}_{d,\gamma,\psi_d(i+1)}$$
 for all $i \in \mathbb{N}$.

Given $d \in \mathbb{N}$, $\tau \in (0,1)$, as well as $c_1 > 0$, and $q_1 \ge 0$ we estimate

$$\sum_{i=\lceil c_1 d^{q_1} \rceil}^{\infty} (\lambda_{d,i})^{\tau} = \sum_{\boldsymbol{m} \in \mathbb{N}^d} (\widetilde{\lambda}_{d,\gamma,\boldsymbol{m}})^{\tau} - \sum_{i=1}^{\lceil c_1 d^{q_1} \rceil - 1} (\lambda_{d,i})^{\tau}$$

$$\geq \prod_{k=1}^d \sum_{m \in \mathbb{N}} (\lambda_{1,\gamma_{d,k},m})^{\tau} - (\lambda_{d,1})^{\tau} (\lceil c_1 d^{q_1} \rceil - 1)$$

$$\geq \prod_{k=1}^d \left(1 + \sum_{k=1}^\infty \left(\frac{\gamma_{d,k}}{\gamma_{d,k} + \pi^2(m-1)^2} \right)^{\tau} \right) - c_1 d^{q_1},$$

since $\lambda_{d,1} = \widetilde{\lambda}_{d,\gamma,(1,\dots,1)} = \prod_{k=1}^d \lambda_{d,\gamma_{d,k},1} = 1$. Due to the boundedness of the generator weights $\gamma_{d,k} \leq C_{\gamma}$ for every $k \in \{1,\dots,d\}$, we can further estimate the sum by

$$\sum_{m=2}^{\infty} \left(\frac{\gamma_{d,k}}{\gamma_{d,k} + \pi^2 (m-1)^2} \right)^{\tau} \geq \gamma_{d,k}^{\tau} \sum_{i=1}^{\infty} \frac{(C_{\gamma}')^{\tau}}{i^{2\tau}} = \gamma_{d,k}^{\tau} \left(C_{\gamma}' \right)^{\tau} \zeta(2\tau),$$

where we set $C'_{\gamma} = (C_{\gamma} + \pi^2)^{-1}$. Because $\ln(1+y) \ge y/(1+y)$ for all $y \ge 0$ we conclude that for $k = 1, \ldots, d$ and some positive C depending on C_{γ} and τ we have

$$\ln(1 + \gamma_{d,k}^{\tau} (C_{\gamma}')^{\tau} \zeta(2\tau)) \ge \frac{(C_{\gamma}')^{\tau} \zeta(2\tau)}{1 + \gamma_{d,k}^{\tau} (C_{\gamma}')^{\tau} \zeta(2\tau)} \gamma_{d,k}^{\tau} \ge C \gamma_{d,k}^{\tau}.$$

Consequently,

$$\sum_{i=\lceil c_1 d^{q_1} \rceil}^{\infty} (\lambda_{d,i})^{\tau} \ge \prod_{k=1}^{d} \exp(C\gamma_{d,k}^{\tau}) - c_1 d^{q_1} = \exp\left(C\sum_{k=1}^{d} \gamma_{d,k}^{\tau}\right) - c_1 d^{q_1}.$$

Therefore, polynomial tractability implies $q(\gamma) < 1$, and strong polynomial tractability is possible only if $p(\gamma) < 1$. Here p and q describe the sum exponents of the product weight sequence $\gamma = (\gamma_{\alpha})_{\alpha \in \mathbb{N}_{d}^{d}}, d \in \mathbb{N}$, defined at the beginning of Section 4.3.2.

Together with Proposition 4.7 this finally proves

THEOREM 4.18. Consider the uniform approximation problem defined on the sequence $(\mathcal{H}_d^{\gamma})_{d\in\mathbb{N}}$ of unanchored Sobolev spaces, where the product weight sequence γ is constructed from a uniformly bounded generator sequence $C_{\gamma} \geq \gamma_{d,1} \geq \cdots \geq \gamma_{d,d}$, $d \in \mathbb{N}$. We study this problem in the worst case setting and with respect to the absolute error criterion. Then we have

- polynomial tractability if and only if $q(\gamma) < 1$, and
- strong polynomial tractability if and only if $p(\gamma) < 1$.

5. Problems on Hilbert spaces with (anti)symmetry conditions

In this last chapter we describe an essentially new kind of a priori knowledge which can help to overcome the curse of dimensionality. As in Section 2.4, we study compact linear problems $S = (S_d)_{d \in \mathbb{N}}$ defined between tensor products of Hilbert spaces but now we restrict our attention to problem elements which fulfill certain (anti)symmetry conditions. After investigating some basic properties of the related subspaces of (anti)symmetric problem elements in Section 5.1 we construct a linear algorithm that uses finitely many continuous linear functionals and show an explicit formula for its worst case error in terms of the eigenvalues $\lambda = (\lambda_m)_{m \in \mathbb{N}}$ of the operator $W_1 = S_1^{\dagger} S_1$. Moreover, in Section 5.2 we show that this algorithm is optimal with respect to a wide class of algorithms. Next we clarify the influence of different (anti)symmetry conditions on the complexity, compared to the case for the classical unrestricted problem studied in Section 2.4.3. In particular, in Section 5.3 we give necessary and sufficient conditions for (strong) polynomial tractability of (anti)symmetric problems. Apart from the absolute error criterion we deal with normalized errors as well. Finally, in Section 5.4, we discuss several applications. Section 5.4.2 in particular indicates how to apply our results to the approximation problem for wavefunctions. For applications of symmetry conditions to integration problems we refer to [Wei14].

Most of the results stated in this chapter are already published in [Wei11] and [Wei12a]. At some points we improve the known results and/or proof techniques slightly. In particular, the results presented also hold for problems defined on finite-dimensional or on non-separable source spaces.

- **5.1.** Basic definitions related to (anti)symmetry. The aim of this section is to introduce the notion of (anti)symmetry in Hilbert spaces. In order to illustrate this concept we mainly deal with *function spaces*. For this purpose in Section 5.1.1 we start by defining (anti)symmetry properties for functions which will lead us to orthogonal projections, mapping the whole space onto its subspace of (anti)symmetric functions. In Section 5.1.2 it will turn out that these projections applied to a given basis of a tensor product Hilbert function space lead us to handsome formulas for orthonormal bases of the subspaces. Finally we generalize our approach and define (anti)symmetry conditions for *arbitrary* tensor product Hilbert spaces based on the results deduced for function spaces. Section 5.1.3 is devoted to this generalization.
- **5.1.1.** Hilbert function spaces. Following Hamaekers [Ham09, Section 2.5] we use a general approach to (anti)symmetric functions which can also be found in [Wei12a].

Let H be a (possibly non-separable) Hilbert space of real-valued multivariate functions f defined on some domain Ω in \mathbb{R}^d , where we assume $d \geq 2$ to be fixed. Furthermore, take an arbitrary non-empty subset of coordinates $I \subseteq \{1, \ldots, d\}$. For every such subset we define the set

$$S_I = \left\{ \pi \colon \{1, \dots, d\} \to \{1, \dots, d\} \mid \pi \text{ bijective and } \pi|_{\{1, \dots, d\} \setminus I} = \mathrm{id} \right\}$$
 (5.1)

of all permutations on $\{1, \ldots, d\}$ that leave the complement of I fixed. To abbreviate the notation we identify $\pi \in \mathcal{S}_I$ with the corresponding permutation π' on \mathbb{R}^d ,

$$\pi' : \mathbb{R}^d \to \mathbb{R}^d, \quad \boldsymbol{x} = (x_1, \dots, x_d) \mapsto \pi'(\boldsymbol{x}) = (x_{\pi(1)}, \dots, x_{\pi(d)}).$$

For an appropriate definition of partial (anti)symmetry of functions $f \in H$ we need the following simple assumptions. Given any $\pi \in \mathcal{S}_I$ we assume that

(A5.1) $\boldsymbol{x} \in \Omega$ implies $\pi(\boldsymbol{x}) \in \Omega$,

(A5.2) $f \in H$ implies $f(\pi(\cdot)) \in H$, and

(A5.3) there exists $c_{\pi} \geq 0$ (independent of f) such that $||f(\pi(\cdot))|| H|| \leq c_{\pi} ||f|| H||$.

A function $f \in H$ is called partially symmetric with respect to I (or I-symmetric for short) if any permutation $\pi \in S_I$ applied to the argument x does not affect the value of f. Hence,

$$f(\mathbf{x}) = f(\pi(\mathbf{x}))$$
 for all $\mathbf{x} \in \Omega$ and every $\pi \in \mathcal{S}_I$. (5.2)

Moreover, we call a function $f \in H$ partially antisymmetric with respect to I (or I-antisymmetric) if f changes its sign on exchanging the variables x_i and x_j with each other, where $i, j \in I$. That is,

$$f(\boldsymbol{x}) = (-1)^{|\pi|} f(\pi(\boldsymbol{x}))$$
 for all $\boldsymbol{x} \in \Omega$ and every $\pi \in \mathcal{S}_I$, (5.3)

where $|\pi|$ denotes the *inversion number* of the permutation π . The term $(-1)^{|\pi|}$ therefore coincides with the *sign*, or *parity*, of π and is equal to the determinant of the associated permutation matrix. In the case #I = 1 we do not claim any (anti)symmetry, since then the set $S_I = \{id\}$ is trivial. For $I = \{1, \ldots, d\}$ functions f which satisfy (5.2) or (5.3), respectively, are called *fully (anti)symmetric*.

Note that, in particular, formula (5.3) implies that the value f(x) of I-antisymmetric functions f equals zero if $x_i = x_j$ with $i \neq j$ and $i, j \in I$. For I-symmetric functions such an implication does not hold. Therefore the (partial) antisymmetry property is a somewhat more restrictive condition than the (partial) symmetry property with respect to the same subset I. As we will see in Section 5.3 this will affect our complexity estimates as well.

Next we define the so-called symmetrizer \mathfrak{S}_I^H and antisymmetrizer \mathfrak{A}_I^H on H with respect to the subset I by

$$\begin{split} \mathfrak{S}_I^H \colon H \to H, \quad f \mapsto \mathfrak{S}_I^H(f) &= \frac{1}{\# \mathcal{S}_I} \sum_{\pi \in \mathcal{S}_I} f(\pi(\boldsymbol{\cdot})), \\ \mathfrak{A}_I^H \colon H \to H, \quad f \mapsto \mathfrak{A}_I^H(f) &= \frac{1}{\# \mathcal{S}_I} \sum_{\pi \in \mathcal{S}_I} (-1)^{|\pi|} f(\pi(\boldsymbol{\cdot})). \end{split}$$

If there is no danger of confusion we use the notation \mathfrak{S}_I and \mathfrak{A}_I instead of \mathfrak{S}_I^H and \mathfrak{A}_I^H , respectively. The following lemma collects some basic properties. It generalizes Lemma 10.1 in Zeiser [Zei10].

LEMMA 5.1. For $\emptyset \neq I \subseteq \{1, ..., d\}$ both the mappings $P_I \in \{\mathfrak{S}_I, \mathfrak{A}_I\}$ define bounded linear operators on the Hilbert space H with $P_I^2 = P_I$. Thus, \mathfrak{S}_I and \mathfrak{A}_I provide projections of H onto the closed linear subspaces

$$\mathfrak{S}_I(H) = \{ f \in H \mid f \text{ satisfies (5.2)} \}$$
 and $\mathfrak{A}_I(H) = \{ f \in H \mid f \text{ satisfies (5.3)} \}$ (5.4)

of all partially (anti)symmetric functions with respect to I in H, respectively. If, in addition,

$$\langle f(\pi(\cdot)), g(\pi(\cdot)) \rangle_H = \langle f, g \rangle_H \quad \text{for all } f, g \in H \text{ and every } \pi \in \mathcal{S}_I$$
 (5.5)

then the operators are self-adjoint and hence the projections are orthogonal. Consequently,

$$H = \mathfrak{S}_I(H) \oplus (\mathfrak{S}_I(H))^{\perp} = \mathfrak{A}_I(H) \oplus (\mathfrak{A}_I(H))^{\perp}. \tag{5.6}$$

Proof. Obviously $P_I \in \{\mathfrak{S}_I, \mathfrak{A}_I\}$ is well-defined due to the assumptions (A5.1) and (A5.2). Linearity follows directly from the definition and, using (A5.3), we see that the operator norm of P_I is bounded by $\max \{c_{\pi} \mid \pi \in \mathcal{S}_I\}$.

To show that the operators are idempotent, i.e. that $P_I^2 = P_I$, we first prove that $\mathfrak{A}_I(f)$ satisfies (5.3) for every $f \in H$. Therefore, we use the representation

$$(\mathfrak{A}_{I}(f))(\pi(\cdot)) = \frac{1}{\#\mathcal{S}_{I}} \sum_{\sigma \in \mathcal{S}_{I}} (-1)^{|\sigma|} f(\sigma(\pi(\cdot))) = \frac{1}{\#\mathcal{S}_{I}} \sum_{\lambda \in \mathcal{S}_{I}} (-1)^{|\lambda| + |\pi|} f(\lambda(\cdot))$$
$$= (-1)^{|\pi|} (\mathfrak{A}_{I}(f))(\cdot)$$

for every fixed $\pi \in \mathcal{S}_I$. Here we imposed $\lambda = \sigma \circ \pi \in \mathcal{S}_I$ and used the fact that

$$|\lambda \circ \pi^{-1}| = |\lambda| + |\pi^{-1}| = |\lambda| + |\pi|.$$

Hence we have shown $\mathfrak{A}_I(H) \subseteq \{f \in H \mid f \text{ satisfies } (5.3)\}$. In a second step, it is easy to check that for every function $g \in H$ which satisfies (5.3) we have $\mathfrak{A}_I(g) = g$. Thus, $\{f \in H \mid f \text{ satisfies } (5.3)\} \subseteq \mathfrak{A}_I(H)$ and \mathfrak{A}_I is a projector onto $\mathfrak{A}_I(H)$. Since the same arguments also apply to the symmetrizer \mathfrak{S}_I , this shows (5.4), as well as $P_I^2 = P_I$ for $P_I \in \{\mathfrak{S}_I, \mathfrak{A}_I\}$.

To prove the self-adjointness of P_I we need to show that for f and g in H we have $\langle P_I f, g \rangle_H = \langle f, P_I g \rangle_H$. To this end, note that (5.5) is equivalent to

$$\langle f(\pi(\cdot)), g \rangle_H = \langle f, g(\sigma(\cdot)) \rangle_H, \quad f, g \in H, \pi \in \mathcal{S}_I,$$

where we set $\sigma = \pi^{-1}$ and used (A5.2). Now the assertion claimed follows from the bilinearity of the inner product $\langle \cdot, \cdot \rangle_H$. Moreover, orthogonality and the decompositions stated in (5.6) are simple consequences.

We note in passing that (5.5) already implies (A5.3). Furthermore, the notion of partially (anti)symmetric functions can be easily extended to more than one subset I. Indeed, consider two non-empty subsets of coordinates $I, J \subset \{1, \ldots, d\}$ with $I \cap J = \emptyset$. Then we call a function $f \in H$ multiply partially (anti)symmetric with respect to I and

J if f satisfies (5.2), or (5.3), respectively, for I and J. Since I and J are disjoint, we observe that $\pi \circ \sigma = \sigma \circ \pi$ for all $\pi \in \mathcal{S}_I$ and $\sigma \in \mathcal{S}_J$. Hence the linear projections $P_I \in \{\mathfrak{S}_I, \mathfrak{A}_I\}$ and $P_J \in \{\mathfrak{S}_J, \mathfrak{A}_J\}$ commute on H. That is, we have $P_I \circ P_J = P_J \circ P_I$. Further extensions to more than two disjoint subsets of coordinates are possible. We will restrict ourselves to the case of at most two coordinate subsets, because in particular wavefunctions can be modeled as functions which are antisymmetric with respect to I and $J = I^c$, where I^c denotes the complement of I in $\{1, \ldots, d\}$; see, e.g., Section 5.4.2.

5.1.2. Tensor products of Hilbert function spaces. In the previous subsection the function space H was a somewhat abstract Hilbert space of d-variate real-valued functions. Indeed, for the definition of (anti)symmetry we do not need to claim any product structure. On the other hand, it is motivated by applications to consider tensor product function spaces; see, e.g., Section 3.6 in Yserentant [Yse10]. In detail, it is well-known that so-called spaces of dominated mixed smoothness, e.g. $W_2^{(1,\dots,1)}(\mathbb{R}^{3d})$, can be represented as certain tensor products; see Section 1.4.2 in Hansen [Han10].

Anyway, let us take into account such a structure, i.e. let us assume that

$$H = H_d = H_1 \otimes \cdots \otimes H_1 \quad (d \geq 2 \text{ times}),$$

where H_1 is a suitable Hilbert space of functions $f: D \to \mathbb{R}$; see also the constructions given in Section 2.4.1. There it is stated that we can construct an orthonormal basis E_d of H_d from a given ONB E_1 of H_1 ; see (2.13). Since now we deal with function spaces, the d-fold simple tensors in E_d are d-variate functions $e_{d,j}: D^d \to \mathbb{R}$. More precisely, they are given by

$$e_{d,\boldsymbol{j}}(\boldsymbol{x}) = \prod_{l=1}^d e_{j_l}(x_l), \quad \text{ where } \boldsymbol{x} = (x_1,\ldots,x_d) \in D^d \text{ and } \boldsymbol{j} \in \mathcal{I}_d = (\mathcal{I}_1)^d,$$

provided that $E_1 = \{e_m : D \to \mathbb{R} \mid m \in \mathcal{I}_1\}$ denotes the underlying ONB in H_1 . To exploit this representation we start with a simple observation.

Let $d \in \mathbb{N}$. Moreover assume $j \in \mathcal{I}_d$ and $x \in D^d$, as well as a non-empty subset I of $\{1, \ldots, d\}$, to be arbitrarily fixed. If we define $\sigma = \pi^{-1} \in \mathcal{S}_I$ then

$$e_{d,j}(\pi(\mathbf{x})) = \prod_{l=1}^{d} e_{j_l}(x_{\pi(l)}) = \prod_{l=1}^{d} e_{j_{\sigma(l)}}(x_l) = e_{d,\sigma(j)}(\mathbf{x}).$$
 (5.7)

For simplicity, once again we identified $\pi(\mathbf{j}) = \pi(j_1, \dots, j_d)$ with $(j_{\pi(1)}, \dots, j_{\pi(d)})$ for $\mathbf{j} \in \mathcal{I}_d = (\mathcal{I}_1)^d$. Since $\mathbf{x} \in D^d$ was arbitrary and $|\pi| = |\pi^{-1}| = |\sigma|$ we obtain

$$\mathfrak{S}_{I}e_{d,j} = \frac{1}{\#\mathcal{S}_{I}} \sum_{\sigma \in \mathcal{S}_{I}} e_{d,\sigma(j)} \quad \text{and} \quad \mathfrak{A}_{I}e_{d,j} = \frac{1}{\#\mathcal{S}_{I}} \sum_{\sigma \in \mathcal{S}_{I}} (-1)^{|\sigma|} e_{d,\sigma(j)}$$
 (5.8)

for all $j \in \mathcal{I}_d$. Besides this, (5.7) can be used to verify that (5.5) in Lemma 5.1 always holds true for (unweighted) tensor products of Hilbert function spaces.

Note that in general, i.e. for arbitrary $j \in \mathcal{I}_d$ and $\sigma \in \mathcal{S}_I$, the tensor products $e_{d,\sigma(j)}$ and $e_{d,j}$ do not coincide, because taking the tensor product is not commutative in general. Therefore \mathfrak{S}_I is not simply the identity on the set of basis functions $E_d = \{e_{d,j} \mid j \in \mathcal{I}_d\}$. On the other hand, we see that for different $j \in \mathcal{I}_d$ many of the functions $\mathfrak{S}_I e_{d,j}$ coincide. Of course the same holds true for $\mathfrak{A}_I e_{d,j}$, at least up to a factor of -1.

We will see in the following that for $P_I \in \{\mathfrak{S}_I, \mathfrak{A}_I\}$ a linearly independent subset of all projections $\{P_I e_{d,j} \mid j \in \mathcal{I}_d\}$ equipped with suitable normalizing constants can be used as an ONB of the linear subspace $P_I(H_d)$ of I-(anti)symmetric functions in H_d . For the application we have in mind, we need this result only in the case where the underlying space H_1 is separable. Without loss of generality, we can thus assume that (2^9)

$$\mathcal{I}_1 = \mathcal{M}_1 = \{ m \in \mathbb{N} \mid m < \dim H_1 + 1 \}$$

and consequently $\mathcal{I}_d = \mathcal{M}_d = (\mathcal{M}_1)^d \subseteq \mathbb{N}^d$. Clearly, in the most interesting case the set \mathcal{I}_d equals \mathbb{N}^d .

To formalize the assertion, we need a further definition. For fixed $d \geq 2$ and $I \subseteq \{1, \ldots, d\}$, let us introduce a function

$$M_I = M_{I,d} \colon \mathbb{N}^d \to \{0, \dots, \#I\}^{\#I}$$

which counts how often different indices occur in a given multi-index $j \in \mathbb{N}^d$ among the subset I of coordinates, ordered with respect to their rate. To give an example let d = 7 and $I = \{1, \ldots, 6\}$. Then $M_{I,7}$ applied to $j = (12, 4, 4, 12, 6, 4, 4) \in \mathbb{N}^7$ gives the (#I = 6)-dimensional vector $M_{I,7}(j) = (3, 2, 1, 0, 0, 0)$, because j contains the number "4" three times among the coordinates j_1, \ldots, j_6 , "12" two times, and so on. Since in this example there are only three different numbers involved, the fourth to sixth coordinates of $M_{I,7}(j)$ equal zero. Obviously, M_I is invariant under all permutations $\pi \in \mathcal{S}_I$ of the argument. Thus,

$$oldsymbol{M}_I(oldsymbol{j}) = oldsymbol{M}_I(\pi(oldsymbol{j})) \quad ext{ for all } oldsymbol{j} \in \mathbb{N}^d ext{ and } \pi \in \mathcal{S}_I.$$

In addition, since $M_I(j)$ is again a multi-index, we see that $|M_I(j)| = \#I$ and $M_I(j)!$ are well-defined for every $j \in \mathbb{N}^d$. With this tool prepared, we are ready to prove

LEMMA 5.2. Assume $E_d = \{e_{d,j} \mid j \in \mathcal{M}_d\}$ is a given orthonormal tensor product basis in the space H_d and let $\emptyset \neq I = \{i_1, \dots, i_{\#I}\} \subseteq \{1, \dots, d\}$. Moreover, for $P_I \in \{\mathfrak{S}_I, \mathfrak{A}_I\}$ define the functions $\xi_j \colon D^d \to \mathbb{R}$ by

$$\xi_{\boldsymbol{j}} = \sqrt{\frac{\#\mathcal{S}_I}{\boldsymbol{M}_I(\boldsymbol{j})!}} P_I(e_{d,\boldsymbol{j}}) \quad \text{ for } \boldsymbol{j} \in \mathcal{M}_d.$$

Then the set $\Xi_d = \{\xi_k \mid k \in \nabla_d\}$ is an orthonormal basis of the partially (anti)symmetric subspace $P_I(H_d)$, where ∇_d is given by

$$\nabla_{d} = \begin{cases} \{ \boldsymbol{k} \in \mathcal{M}_{d} \mid k_{i_{1}} \leq \dots \leq k_{i_{\#I}} \} & \text{if } P_{I} = \mathfrak{S}_{I}, \\ \{ \boldsymbol{k} \in \mathcal{M}_{d} \mid k_{i_{1}} < \dots < k_{i_{\#I}} \} & \text{if } P_{I} = \mathfrak{A}_{I}. \end{cases}$$

$$(5.9)$$

Proof. To abbreviate the notation, we suppress the index H_d on the inner products $\langle \cdot, \cdot \rangle_{H_d}$ in this proof.

 $^(^{29})$ Note that also the case of abstract, countable index sets \mathcal{I}_1 can be reduced to this form by the application of some simple isomorphism.

Step 1. We start by proving orthonormality. Therefore let us recall (5.8) and remember that now $\mathcal{I}_d = \mathcal{M}_d$. For $P_I = \mathfrak{A}_I$ and $j, k \in \nabla_d$ easy calculations yield

$$\langle \xi_{\boldsymbol{j}}, \xi_{\boldsymbol{k}} \rangle = \frac{\# \mathcal{S}_{I}}{\sqrt{\boldsymbol{M}_{I}(\boldsymbol{j})! \cdot \boldsymbol{M}_{I}(\boldsymbol{k})!}} \langle \mathfrak{A}_{I}(e_{d,\boldsymbol{j}}), \mathfrak{A}_{I}(e_{d,\boldsymbol{k}}) \rangle$$

$$= \frac{1}{\# \mathcal{S}_{I} \sqrt{\boldsymbol{M}_{I}(\boldsymbol{j})! \cdot \boldsymbol{M}_{I}(\boldsymbol{k})!}} \sum_{\pi, \sigma \in \mathcal{S}_{I}} (-1)^{|\pi| + |\sigma|} \langle e_{d,\pi(\boldsymbol{j})}, e_{d,\sigma(\boldsymbol{k})} \rangle.$$

Of course, up to the factor controlling the sign, the same is true for the case $P_I = \mathfrak{S}_I$. Now assume that there exists $l \in \{1, \ldots, d\}$ such that $j_l \neq k_l$. Then the ordering of $j, k \in \nabla_d$ implies that $\pi(j) \neq \sigma(k)$ for all $\sigma, \pi \in \mathcal{S}_I$, since π and σ leave the coordinates $l \in I^c$ fixed. Hence, we conclude that $\pi(j) = \sigma(k)$ only if j = k.

At this point we have to distinguish the antisymmetric and the symmetric cases. For $P_I = \mathfrak{A}_I$ the only way to deduce $\pi(j) = \sigma(k)$ is to assume j = k and $\pi = \sigma$. Furthermore we see that in the antisymmetric case we have $M_I(j)! = 1$ for all $j \in \nabla_d$, because then all coordinates j_l , where $l \in I$, differ. Therefore, in this case the last inner product coincides with $\delta_{j,k}\delta_{\pi,\sigma}$ because of the mutual orthonormality of the elements from $E_d = \{e_{d,j} \mid j \in \mathcal{M}_d\}$. Hence we arrive at

$$\langle \xi_{\boldsymbol{j}}, \xi_{\boldsymbol{k}} \rangle = \frac{1}{\# \mathcal{S}_I} \sum_{\pi \in \mathcal{S}_I} (-1)^{2|\pi|} \delta_{\boldsymbol{j}, \boldsymbol{k}} = \delta_{\boldsymbol{j}, \boldsymbol{k}} \quad \text{ for all } \boldsymbol{j}, \boldsymbol{k} \in \nabla_d,$$

as claimed.

So, let us consider the case $P_I = \mathfrak{S}_I$ and $j = k \in \nabla_d$, since we already saw that otherwise $\langle \xi_j, \xi_k \rangle$ equals zero. Then for fixed $\sigma \in \mathcal{S}_I$ there are $M_I(j)$! different permutations $\pi \in \mathcal{S}_I$ such that $\pi(j) = \sigma(j)$. This leads to

$$\langle \xi_{j}, \xi_{j} \rangle = \frac{1}{\# \mathcal{S}_{I} \cdot M_{I}(j)!} \sum_{\sigma \in \mathcal{S}_{I}} M_{I}(j)! = 1$$

and completes the proof of orthonormality.

Step 2. It remains to show that the span of $\Xi_d = \{\xi_k \mid k \in \nabla_d\}$ is dense in $P_I(H_d)$ for $P_I \in \{\mathfrak{S}_I, \mathfrak{A}_I\}$. Note that every multi-index $j \in \mathcal{M}_d$ can be represented by a uniquely defined multi-index $k \in \nabla_d$ and exactly $M_I(k)$! different permutations $\pi \in \mathcal{S}_I$ such that $j = \pi(k)$. Assume that $f \in \mathfrak{A}_I(H_d)$, i.e. $f \in H_d$ satisfies (5.3). Then (5.7) together with (5.5) yields

$$\langle f, e_{d, \mathbf{j}} \rangle = (-1)^{|\pi|} \langle f, e_{d, \pi(\mathbf{j})} \rangle$$
 for all $\mathbf{j} \in \mathcal{M}_d$ and $\pi \in \mathcal{S}_I$. (5.10)

Now expanding f with respect to the basis functions in $E_d \subset H_d$ gives

$$f = \sum_{\boldsymbol{j} \in \mathcal{M}_d} \langle f, e_{d, \boldsymbol{j}} \rangle e_{d, \boldsymbol{j}} = \sum_{\boldsymbol{k} \in \nabla_d} \sum_{\pi \in \mathcal{S}_I} \frac{\langle f, e_{d, \pi(\boldsymbol{k})} \rangle e_{d, \pi(\boldsymbol{k})}}{\boldsymbol{M}_I(\boldsymbol{k})!}$$

$$= \sum_{\boldsymbol{k} \in \nabla_d} \frac{1}{\boldsymbol{M}_I(\boldsymbol{k})!} \sum_{\pi \in \mathcal{S}_I} (-1)^{|\pi|} \langle f, e_{d, \boldsymbol{k}} \rangle e_{d, \pi(\boldsymbol{k})}$$

$$= \sum_{\boldsymbol{k} \in \nabla_d} \sqrt{\frac{\#\mathcal{S}_I}{\boldsymbol{M}_I(\boldsymbol{k})!}} \langle f, e_{d, \boldsymbol{k}} \rangle \sqrt{\frac{\#\mathcal{S}_I}{\boldsymbol{M}_I(\boldsymbol{k})!}} \mathfrak{A}_I(e_{d, \boldsymbol{k}}),$$

where we used (5.8) for the last equality. Furthermore, due to the self-adjointness of \mathfrak{A}_I , we have $\langle f, e_{d,\mathbf{k}} \rangle = \langle \mathfrak{A}_I f, e_{d,\mathbf{k}} \rangle = \langle f, \mathfrak{A}_I e_{d,\mathbf{k}} \rangle$, so that finally $f \in \mathfrak{A}_I(H_d)$ possesses the representation

$$f = \sum_{\mathbf{k} \in \nabla_d} \langle f, \xi_{\mathbf{k}} \rangle \, \xi_{\mathbf{k}}$$

since $\xi_{\mathbf{k}} = \sqrt{\#\mathcal{S}_I/M_I(\mathbf{k})!} \, \mathfrak{A}_I(e_{d,\mathbf{k}})$ per definition. This proves the assertion for $P_I = \mathfrak{A}_I$. The remaining case $P_I = \mathfrak{S}_I$ can be treated in the same way.

Observe that in the antisymmetric case the definition of ξ_{j} for $j \in \nabla_{d}$ simplifies, since then $M_{I}(j)! = 1$ for all $j \in \nabla_{d}$. Moreover we see that in this case ∇_{d} is trivial if $d > \# \mathcal{M}_{1}$. Hence we should assume that dim H_{1} is infinite in order to work with antisymmetric tensor products for arbitrarily many building blocks. We note in passing that the square of the normalizing factor, $\# \mathcal{S}_{I}/M_{I}(j)!$, coincides with the multinomial coefficient $\binom{|M_{I}(j)|}{M_{I}(j)}$ which is quite natural due to combinatorial issues. Furthermore, in the special case $I = \{1, \ldots, \# I\}$ we have

$$P_I(H_d) = P_I\left(\bigotimes_{m \in I} H_1\right) \otimes \left(\bigotimes_{m \notin I} H_1\right).$$

That is, we can consider the subspace of I-(anti)symmetric functions $f \in H_d$ as the tensor product of the set of all fully (anti)symmetric #I-variate functions with the (d - #I)-fold tensor product of H_1 . If #I = 1, i.e. if we do not claim any (anti)symmetry, then $P_I(H_d) = H_d$ and thus we have $\nabla_d = \mathcal{M}_d$, as well as $\Xi_d = E_d$. Modifications in connection with multiple partially (anti)symmetric functions are obvious.

5.1.3. Arbitrary tensor product Hilbert spaces. Up to now we exclusively dealt with Hilbert *function* spaces. However, the proofs of Lemmas 5.1 and 5.2 yield that there are only a few key arguments in connection with (anti)symmetry such that we can cut out this restriction. We briefly sketch the points which need to be changed.

Starting from the very beginning we have to adapt the definition of I-(anti)symmetry due to (5.2) and (5.3) in Section 5.1.1. Of course it is sufficient to define this property at first only for basis elements. Therefore, if $E_d = \{e_{d,\mathbf{k}} \mid \mathbf{k} \in (\mathcal{I}_1)^d = \mathcal{I}_d\}$ denotes a tensor product ONB of H_d and $\emptyset \neq I \subseteq \{1,\ldots,d\}$ is given then we call an element $e_{d,\mathbf{k}} = \bigotimes_{l=1}^d e_{k_l}$ partially symmetric with respect to I (briefly I-symmetric) if

$$e_{d,\mathbf{k}} = e_{d,\pi(\mathbf{k})}$$
 for all $\pi \in \mathcal{S}_I$,

where S_I and $\pi(\mathbf{k}) = (k_{\pi(1)}, \dots, k_{\pi(d)}) \in \mathcal{I}_d$ are defined as before. Analogously, we define I-antisymmetry with an additional factor $(-1)^{|\pi|}$. Finally, an arbitrary element in H_d is called I-(anti)symmetric if in its basis expansion every element with non-vanishing coefficient possesses this property (30).

Next, the antisymmetrizer \mathfrak{A}_I is given as the uniquely defined continuous extension of the linear mapping

$$\widetilde{\mathfrak{A}}_I \colon E_d \to H_d, \quad e_{d,\mathbf{k}} \mapsto \frac{1}{\#\mathcal{S}_I} \sum_{\pi \in \mathcal{S}_I} (-1)^{|\pi|} e_{d,\pi(\mathbf{k})},$$
 (5.11)

⁽³⁰⁾ Even in the non-separable case any such expansion only has countably many terms.

from E_d to H_d . Again the *symmetrizer* \mathfrak{S}_I is given in a similar way. Hence, in the general setting we define the mappings using formula (5.8) which we derived for the special case of function spaces. Note that the triangle inequality yields $||P_I|| \leq 1$ for $P_I \in \{\mathfrak{S}_I, \mathfrak{A}_I\}$.

Once more we denote the sets of all I-(anti)symmetric elements of H_d by $P_I(H_d)$, where $P_I \in \{\mathfrak{S}_I, \mathfrak{A}_I\}$. Observe that this can be justified since the operators P_I again provide orthogonal projections onto closed linear subspaces. That is, a generalization of Lemma 5.1 remains valid also in the more general case of tensor products of arbitrary Hilbert spaces which we consider here. This can be shown using (5.11) and its analogue for \mathfrak{S}_I , as well as with the help of some simple extension arguments. Moreover, also the proof of Lemma 5.2 can be adapted to the generalized setting. Indeed, the only difference is the conclusion of formula (5.10) in Step 2. Now, for arbitrary Hilbert spaces, this simply follows from our definitions. Then the rest of the proof transfers literally.

Finally, without going into details, we stress the point that further generalizations are possible. Here we can think of tensor products of arbitrary Hilbert spaces with multiple partial (anti)symmetry conditions or of scaled tensor products in the sense of Chapter 3. Since the corresponding calculations are straightforward we leave them to the reader.

5.2. Optimal algorithms for (anti)symmetric problems. Keeping the definitions and assertions from Section 5.1 in mind, we are ready to study algorithms for linear problems defined on (anti)symmetric subsets of tensor product Hilbert spaces.

Let $S_d \colon H_d \to \mathcal{G}_d$ denote a tensor product problem in the sense of Section 2.4. It is constructed from a compact linear operator $S_1 \colon H_1 \to \mathcal{G}_1$ between arbitrary Hilbert spaces H_1 and \mathcal{G}_1 via a tensor product construction; see Section 2.4.1. Hence, let $H_d = H_1 \otimes \cdots \otimes H_1$ in what follows and refer to the problem of approximating $S = (S_d)_{d \in \mathbb{N}}$ as the *entire d-variate problem*. Note that we completely solved this problem in Section 2.4. In detail, the *n*th optimal algorithm $A_{n,d}^*$, given by (2.16), was related to a certain subset $\{e_{d,j} = \widetilde{\phi}_{d,j} \mid j \in \mathcal{M}_d\}$ of a tensor product ONB.

In contrast, now we are interested in the approximation of the restriction

$$S_{d,I_d} = S_d \big|_{P_{I_d}(H_d)} \colon P_{I_d}(H_d) \to \mathcal{G}_d$$

of S_d to some (anti)symmetric subspace $P_{I_d}(H_d)$ as defined in Section 5.1.3, where $P_{I_d} \in \{\mathfrak{S}_{I_d}, \mathfrak{A}_{I_d}\}$ and $\emptyset \neq I_d \subseteq \{1, \ldots, d\}$ for $d \in \mathbb{N}$. We refer to $S_I = (S_{d,I_d})_{d \in \mathbb{N}}$ as the I-(anti)symmetric problem. Using the notation from Section 1.2 we thus have $\mathcal{F}_d = P_{I_d}(H_d)$, and consequently $\widetilde{\mathcal{F}}_d = \mathcal{B}(P_{I_d}(H_d))$.

Due to (5.10) it is quite clear that $A_{n,d}^*$ cannot be optimal in this restricted setting since it calculates redundant pieces of information. Hence we need to go beyond this naive attempt to solve I-(anti)symmetric problems efficiently. On the other hand, $P_{I_d}(H_d)$ equipped with the inner product of H_d , $\langle \cdot, \cdot \rangle_{H_d}$ is again a Hilbert space. Therefore we basically know how to construct an optimal algorithm; see Section 2.3.2. If $\#I_d=1$ then our new algorithm should resemble $A_{n,d}^*$, because then we do not claim any (anti)symmetry and thus we deal with the entire tensor product problem.

Before we state the main assertion of this section we present an auxiliary result which shows that any optimal algorithm A^* for S_{d,I_d} needs to preserve the (anti)symmetry

properties of its domain of definition. The following proposition generalizes Lemma 10.2 in Zeiser [Zei10] where this assertion was shown for the approximation problem, that is, for $S_{d,I_d} = \text{id} : P_{I_d}(H_d) \to \mathcal{G}_d$.

PROPOSITION 5.3. Let d > 1 and $\emptyset \neq I \subseteq \{1, ..., d\}$ be arbitrarily fixed. Furthermore, for $X \in \{H, \mathcal{G}\}$ let P_I^X denote the (anti)symmetrizer $P_I \in \{\mathfrak{S}_I, \mathfrak{A}_I\}$ on X_d with respect to I. Then

$$(S_d \circ P_I^H)(g) = (P_I^G \circ S_d)(g) \quad \text{for any } g \in H_d.$$
 (5.12)

Moreover, for all $A: P_I^H(H_d) \to \mathcal{G}_d$ and every $f \in P_I^H(H_d)$,

$$||S_{d,I}f - Af||\mathcal{G}_d||^2 = ||S_{d,I}f - P_I^{\mathcal{G}}(Af)||\mathcal{G}_d||^2 + ||Af - P_I^{\mathcal{G}}(Af)||\mathcal{G}_d||^2.$$
 (5.13)

Hence an optimal algorithm A^* for $S_{d,I}$ preserves (anti)symmetry, i.e.

$$A^* f \in P_I^{\mathcal{G}}(\mathcal{G}_d)$$
 for all $f \in P_I^H(H_d)$.

Proof. First we show that the tensor product operator S_d and the (anti)symmetrizer P_I commute on H_d , i.e. (5.12) holds. In a second step we deduce (5.13). The (anti)symmetry of A^*f for an optimal algorithm A^* then follows immediately.

Step 1. Assume $E_d = \{e_{d,j} \mid j \in \mathcal{I}_d\}$ to be an arbitrary tensor product ONB of H_d , as defined in (2.13). Then, for fixed $j \in \mathcal{I}_d$, formula (5.11) and the structure of $S_d = S_1 \otimes \cdots \otimes S_1$ yield, in the case $P_I = \mathfrak{A}_I$,

$$S_{d}(\mathfrak{A}_{I}^{H}(e_{d,j})) = S_{d}\left(\frac{1}{\#S_{I}} \sum_{\pi \in S_{I}} (-1)^{|\pi|} \bigotimes_{l=1}^{d} e_{j_{\pi(l)}}\right)$$

$$= \frac{1}{\#S_{I}} \sum_{\pi \in S_{I}} (-1)^{|\pi|} \bigotimes_{l=1}^{d} S_{1}(e_{j_{\pi(l)}}) = \mathfrak{A}_{I}^{\mathcal{G}}(S_{d}(e_{d,j})).$$

Obviously the same is true for $P_I = \mathfrak{S}_I$. Hence, (5.12) holds at least on the set of basis elements E_d of H_d . Because of the representation $g = \sum_{j \in \mathcal{I}_d} \langle g, e_{d,j} \rangle_{H_d} e_{d,j}$ of $g \in H_d$, as well as the linearity and boundedness of the operators P_I^H , $P_I^{\mathcal{G}}$ and S_d , we can extend the relation (5.12) from E_d to the whole space H_d .

Step 2. Now let $f \in P_I^H(H_d)$ and let Af denote an arbitrary approximation to $S_{d,I}f$. Then $S_{d,I}f = S_d(P_I^H f) = P_I^{\mathcal{G}}(S_d f)$, due to Step 1. Using the fact that $P_I^{\mathcal{G}}$ provides an orthogonal projection onto $P_I^{\mathcal{G}}(\mathcal{G}_d)$ (see (5.6)) we obtain (5.13), i.e.

$$||S_{d,I}f - Af||\mathcal{G}_{d}||^{2} = ||P_{I}^{\mathcal{G}}(S_{d}f) - [P_{I}^{\mathcal{G}}(Af) + (\mathrm{id}^{\mathcal{G}} - P_{I}^{\mathcal{G}})(Af)]||\mathcal{G}_{d}||^{2}$$

$$= ||P_{I}^{\mathcal{G}}(S_{d}f - Af)||\mathcal{G}_{d}||^{2} + ||(\mathrm{id}^{\mathcal{G}} - P_{I}^{\mathcal{G}})(Af)|||\mathcal{G}_{d}||^{2}$$

$$= ||S_{d,I}f - P_{I}^{\mathcal{G}}(Af)||\mathcal{G}_{d}||^{2} + ||Af - P_{I}^{\mathcal{G}}(Af)|||\mathcal{G}_{d}||^{2},$$

as claimed. \blacksquare

Apart from this qualitative assertion, we are interested in an explicit formula for the optimal algorithm, as well as in sharp error bounds. To get those, let $d \in \mathbb{N}$ and $\emptyset \neq I_d = \{i_1, \ldots, i_{\#I}\} \subseteq \{1, \ldots, d\}$, as well as $P \in \{\mathfrak{S}, \mathfrak{A}\}$. Furthermore, consider the singular value decomposition of $S_1 \colon H_1 \to \mathcal{G}_1$. That is, let $\{(\lambda_m, \phi_m) \mid m \in \mathcal{M}_1\}$ denote the non-trivial eigenpairs of $W_1 = S_1^{\dagger} S_1$; see Section 2.3.1. Due to Proposition 2.10

in Section 2.4.2, we know that for d>1 the (tensor) product eigenpairs $\{(\widetilde{\lambda}_{d,\boldsymbol{m}},\widetilde{\phi}_{d,\boldsymbol{m}})\mid \boldsymbol{m}\in\mathcal{M}_d\}$ of $W_d=S_d^\dagger S_d$ are given by (2.15). Moreover, $E_d=\Phi_d=\{\widetilde{\phi}_{d,\boldsymbol{m}}\mid \boldsymbol{m}\in\mathcal{M}_d\}$ is a tensor product ONB in H_d . Hence, we can apply Lemma 5.2 to $e_{d,\boldsymbol{j}}=\widetilde{\phi}_{d,\boldsymbol{j}},\,\boldsymbol{j}\in\mathcal{M}_d$, in order to obtain an orthonormal basis $\Xi_d=\{\widetilde{\xi}_{\boldsymbol{k}}\mid \boldsymbol{k}\in\nabla_d\}$ of the partially (anti)symmetric subspaces $P_{I_d}(H_d)$. More precisely, for $\boldsymbol{k}\in\nabla_d$ we define

$$\widetilde{\xi}_{k} = \sqrt{\frac{\#S_{I}}{M_{I}(k)!}} P_{I_{d}}\left(\bigotimes_{l=1}^{d} \phi_{k_{l}}\right) \in P_{I}(H_{d}) \quad \text{and} \quad \widetilde{\lambda}_{d,k} = \prod_{l=1}^{d} \lambda_{k_{l}} > 0,$$
(5.14)

where ∇_d is given by (5.9). Similar to the approach in Section 2.4.2, let

$$\psi = \psi_d \colon \{i \in \mathbb{N} \mid i < \#\nabla_d + 1\} \to \nabla_d$$

denote a bijection which provides a non-increasing ordering of $\{\widetilde{\lambda}_{d,\mathbf{k}} \mid \mathbf{k} \in \nabla_d\}$, and put $\lambda_{d,i} = \widetilde{\lambda}_{d,\psi(i)}$, as well as $\xi_{d,i} = \widetilde{\xi}_{\psi(i)}$ for $i < \#\nabla_d + 1$. Finally, if $\#\nabla_d$ is finite then we extend the sequence of λ 's by setting $\lambda_{d,i} = 0$ for $i > \#\nabla_d$.

Given this bunch of notations we are well-prepared to prove our main theorem of this section. For every $d \in \mathbb{N}$ it provides a linear algorithm $A'_{n,d}$ which uses at most n continuous linear functionals on the input to approximate the solution operator S_{d,I_d} of a given I_d -(anti)symmetric tensor product problem between Hilbert spaces. Since the worst case error of this algorithm coincides with the nth minimal error of the problem, $A'_{n,d}$ is optimal in this setting; thus it cannot be improved by any other algorithm from the class $\mathcal{A}_d^{n,\text{cont}} \cup \mathcal{A}_d^{n,\text{adapt}}$; see Section 1.3. The assertion reads as follows.

THEOREM 5.4. Assume that $S_I = (S_{d,I_d})_{d \in \mathbb{N}}$ is the linear tensor product problem S restricted to the I_d -(anti)symmetric subspaces $P_{I_d}(H_d)$ of the d-fold tensor product spaces H_d . Then for every $d \in \mathbb{N}$,

$$\{(\lambda_{d,i}, \xi_{d,i}) \mid 1 \le i < \#\nabla_d + 1\} = \{(\widetilde{\lambda}_{d,k}, \widetilde{\xi}_k) \mid k \in \nabla_d\}$$

$$(5.15)$$

is the set of eigenpairs of $W_{d,I_d} = S_{d,I_d}^{\dagger} S_{d,I_d} \colon P_{I_d}(H_d) \to P_{I_d}(H_d)$. Thus, for every $n \in \mathbb{N}_0$ and all $d \in \mathbb{N}$, the linear algorithm $A'_{n,d} \colon P_{I_d}(H_d) \to P_{I_d}(\mathcal{G}_d)$ given by

$$A'_{n,d}f = \sum_{i=1}^{\min\{n, \#\nabla_d\}} \langle f, \xi_{d,i} \rangle_{H_d} S_d \xi_{d,i},$$
 (5.16)

is nth optimal for S_{d,I_d} with respect to the worst case setting. Furthermore we have

$$e^{\text{wor}}(n, d; P_{I_d}(H_d)) = \Delta^{\text{wor}}(A'_{n,d}; P_{I_d}(H_d)) = \sqrt{\lambda_{d,n+1}}.$$
 (5.17)

Proof. Since S_I is a compact problem between Hilbert spaces it is enough to prove that for $d \in \mathbb{N}$ the eigenpairs of $W_{d,I_d} = S_{d,I_d}^{\dagger} S_{d,I_d}$ are given by (5.15). The remaining assertions then follow from Corollary 2.7. Indeed, we only need to show that $W_{d,I_d} \widetilde{\xi}_{\boldsymbol{k}} = \widetilde{\lambda}_{d,\boldsymbol{k}} \widetilde{\xi}_{\boldsymbol{k}}$ for every $\boldsymbol{k} \in \nabla_d$ because we already know that the set $\Xi_d = \{\widetilde{\xi}_{\boldsymbol{k}} \mid \boldsymbol{k} \in \nabla_d\}$ builds an ONB in $P_{I_d}(H_d)$. Hence there cannot be more than these eigenpairs.

To prove the claim, observe that from the first part of Proposition 5.3 it follows that

$$S_{d,I_d} = S_d \circ P_{I_d}^H = P_{I_d}^{\mathcal{G}} \circ S_d, \quad \text{which implies} \quad S_{d,I_d} \colon P_{I_d}^H(H_d) \to P_{I_d}^{\mathcal{G}}(\mathcal{G}_d).$$

Moreover, due to the self-adjointness of the projectors (see Lemma 5.1), it is easily seen that this yields

$$S_{d,I_d}^\dagger = P_{I_d}^H \circ S_d^\dagger = S_d^\dagger \circ P_{I_d}^{\mathcal{G}} \quad \text{so that} \quad S_{d,I_d}^\dagger \colon P_{I_d}^{\mathcal{G}}(\mathcal{G}_d) \to P_{I_d}^H(H_d).$$

Consequently, we have

$$W_{d,I_d} P_{I_d}^H = (P_{I_d}^H S_d^{\dagger})(S_d P_{I_d}^H) P_{I_d}^H = P_{I_d}^H (S_d^{\dagger} P_{I_d}^{\mathcal{G}}) S_d = P_{I_d}^H (S_d^{\dagger} S_d) = P_{I_d}^H W_d,$$

because of $(P_{I_d}^X)^2 = P_{I_d}^X$, where $X \in \{H, \mathcal{G}\}$. Since for every $j \in \mathcal{M}_d$ the simple tensor $\widetilde{\phi}_{d,j}$ is an eigenelement of W_d with respect to the eigenvalue $\widetilde{\lambda}_{d,j}$, we conclude that

$$W_{d,I_d}(P_{I_d}^H \widetilde{\phi}_{d,\boldsymbol{j}}) = \widetilde{\lambda}_{d,\boldsymbol{j}}(P_{I_d}^H \widetilde{\phi}_{d,\boldsymbol{j}})$$

from the linearity of $P_{I_d}^H$. In particular, this is true for every $j = k \in \nabla_d \subseteq \mathcal{M}_d$. But now we note that $\widetilde{\xi}_k$ equals $P_{I_d}^H \widetilde{\phi}_{d,k}$, at least up to some normalizing constant. Hence, using linearity once again, we have proven the assertion.

We conclude this section by adding some final remarks on the above theorem.

REMARK 5.5. Obviously, our former result for the entire tensor product problem $S = (S_d : H_d \to \mathcal{G}_d)_{d \in \mathbb{N}}$ in Section 2.4.2 is covered by Theorem 5.4 as well. We simply have to choose I_d such that $\#I_d = 1$ for every $d \in \mathbb{N}$ and obtain $A'_{n,d} = A^*_{n,d}$. As in this case, the worst case error can be attained by the element $\xi_{d,n+1}$ provided that $n < \#\nabla_d$. Otherwise it trivially equals zero.

It should be clear to the reader how to generalize the results of this section to the case of multiple partially (anti)symmetric problems where we claim (anti)symmetry with respect to more than one subset of coordinates I. Recall that this definition is given at the end of Section 5.1.1.

Finally we want to mention that we decided to give a different proof of Theorem 5.4 than those in [Wei11] and [Wei12a]. The reason is that the usage of the self-adjointness of the projections P_{I_d} seems to be more elegant than again repeating the arguments used for Corollary 2.7 in Section 2.3.2. Furthermore, now we can handle also problems defined on non-separable or on finite-dimensional source spaces H_d . Thus we slightly generalized our old results. \Box

5.3. Complexity of (anti)symmetric problems. Encouraged by the exact formula for the nth minimal worst case error in Theorem 5.4 the intention of the present section is to investigate the information complexity of (anti)symmetric tensor product problems. We restrict our attention to the study of polynomial and strong polynomial tractability in what follows. The aim is to find necessary and sufficient conditions for these properties in terms of the univariate sequence $\lambda = (\lambda_m)_{m \in \mathbb{N}}$ and the number of (anti)symmetry conditions we impose. From the definition of ∇_d in (5.9) it is quite clear that antisymmetric problems are significantly easier than their symmetric counterparts. Therefore, after proving some general assertions, we handle these cases separately in order to obtain sharp conditions. Moreover, we distinguish between the absolute and the normalized error criterions.

Let us fix the basic notation for this section. As before, assume $S_I = (S_{d,I_d})_{d \in \mathbb{N}}$ denotes a tensor product problem $S = (S_d \colon H_d \to \mathcal{G}_d)_{d \in \mathbb{N}}$, restricted to some sequence of (anti)symmetric subspaces $P_{I_d}(H_d)$, where $P_{I_d} \in \{\mathfrak{S}_{I_d}, \mathfrak{A}_{I_d}\}$, of the tensor product Hilbert spaces $H_d = H_1 \otimes \cdots \otimes H_d$, $d \in \mathbb{N}$. Here for every $d \in \mathbb{N}$ the elements are (anti)symmetric with respect to the non-empty subset $I_d \subseteq \{1,\ldots,d\}$ of coordinates. The cardinality of those subsets will be denoted by $a_d = \#I_d$ and we set $b_d = d - a_d$ for the number of coordinates without (anti)symmetry conditions. Finally, for $d \in \mathbb{N}$ the non-increasingly ordered eigenvalues $\lambda_{d,i} = \widetilde{\lambda}_{d,\psi(i)}$, $i \in \mathbb{N}$, are given by (5.14) and (5.15). They are constructed out of the squared singular values $\lambda = (\lambda_m)_{m \in \mathbb{N}}$ of the underlying solution operator $S_1 \colon H_1 \to \mathcal{G}_1$.

As an immediate consequence of (5.17) we see that the initial error of approximating S_{d,I_d} on the unit ball $\widetilde{\mathcal{F}}_d = \mathcal{B}(P_{I_d}(H_d))$ is given by

$$\varepsilon_d^{\text{init}} = e^{\text{wor}}(0, d; P_{I_d}(H_d)) = \sqrt{\lambda_{d,1}} = \begin{cases} \sqrt{\lambda_1^d} & \text{if } P_{I_d} = \mathfrak{S}_{I_d}, \\ \sqrt{\lambda_1^{b_d} \cdot \lambda_1 \cdot \ldots \cdot \lambda_{a_d}} & \text{if } P_{I_d} = \mathfrak{A}_{I_d}. \end{cases}$$

Clearly, we need to assume that this initial error is strictly positive for any reasonably large $d \in \mathbb{N}$ because otherwise we have (strong) polynomial tractability by default. In particular, if the number of antisymmetric coordinates a_d grows with the dimension then this condition implies that the whole sequence of univariate eigenvalues λ need to be strictly positive. Moreover, similar to the entire tensor product problems studied in Section 2.4.3, we always assume that $\lambda_2 > 0$ in order to avoid triviality. Consequently, we have $\#\mathcal{M}_1 \geq 2$.

Now we are ready to present a first general condition which is necessary for (strong) polynomial tractability of both symmetric and antisymmetric problems as long as we deal with the absolute error criterion. It is independent of the concrete choice of the (anti)symmetry conditions we impose.

LEMMA 5.6 (General necessary conditions, absolute errors). Let $P_{I_d} \in \{\mathfrak{S}_{I_d}, \mathfrak{A}_{I_d}\}$ and consider $S_I = (S_{d,I_d})_{d \in \mathbb{N}}$ as defined above, where I_d is arbitrarily fixed for every $d \in \mathbb{N}$. Then the fact that S_I is polynomially tractable with the constants C, p > 0 and $q \geq 0$ implies that $\lambda = (\lambda_m)_{m \in \mathbb{N}} \in \ell_{\tau}$ for all $\tau > p/2$. Moreover, for any such τ and all $d \in \mathbb{N}$,

$$\frac{1}{(\lambda_{d,1})^{\tau}} \sum_{\boldsymbol{k} \in \nabla_d} (\widetilde{\lambda}_{d,\boldsymbol{k}})^{\tau} \le (1+C) d^q + C^{2\tau/p} \zeta \left(\frac{2\tau}{p}\right) \left(\frac{d^{2q/p}}{\lambda_{d,1}}\right)^{\tau}.$$

Proof. From Theorem 2.8 we know that for any $\tau > p/2$ and r = 2q/p polynomial tractability yields

$$\sup_{d \in \mathbb{N}} \frac{1}{d^r} \left(\sum_{i=f(d)}^{\infty} (\lambda_{d,i})^{\tau} \right)^{1/\tau} < \infty, \tag{5.18}$$

where the function $f: \mathbb{N} \to \mathbb{N}$ is given by $f(d) = \lceil (1+C) d^q \rceil$. This in particular implies that the sum in brackets converges for every fixed $d \in \mathbb{N}$. Therefore, in particular for d = 1 the tail series $\sum_{i=f(1)}^{\infty} (\lambda_{1,i})^{\tau} = \sum_{m=\lceil 1+C \rceil}^{\infty} (\lambda_m)^{\tau}$ needs to be finite, which is possible only if $\lambda = (\lambda_m)_{m \in \mathbb{N}} \in \ell_{\tau}$.

So, let us turn to the second assertion. Obviously (5.18) implies the existence of some constant $C_1 > 0$ such that

$$\sum_{i=f(d)}^{\infty} (\lambda_{d,i})^{\tau} \le C_1 d^{r\tau} \quad \text{ for all } d \in \mathbb{N}.$$

Indeed, Theorem 2.8 implies that we can take $C_1 = C^{2\tau/p}\zeta(2\tau/p)$. Due to the ordering of $(\lambda_{d,i})_{i\in\mathbb{N}}$ the rest of the sum can also be bounded easily for any $d\in\mathbb{N}$ by

$$\sum_{i=1}^{f(d)-1} (\lambda_{d,i})^{\tau} \le (\lambda_{d,1})^{\tau} (f(d)-1).$$

Since $\sum_{\mathbf{k} \in \nabla_d} (\widetilde{\lambda}_{d,\mathbf{k}})^{\tau} = \sum_{i=1}^{\infty} (\lambda_{d,i})^{\tau}$, it remains to show that $f(d) - 1 \leq (1+C)d^q$ for every $d \in \mathbb{N}$, which is also obvious due to the definition of f.

5.3.1. Symmetric problems (absolute errors). Apart from the general assertion $\lambda \in \ell_{\tau}$, we focus our attention on further necessary conditions for (strong) polynomial tractability in the symmetric setting. The following proposition yields a slight improvement compared to the corresponding assertion stated in [Wei12a] which can be obtained without using essentially new ideas.

PROPOSITION 5.7 (Necessary conditions, symmetric case). Let $S_I = (S_{d,I_d})_{d \in \mathbb{N}}$ be the problem considered in Lemma 5.6 and set $P_{I_d} = \mathfrak{S}_{I_d}$. Moreover, assume $\lambda_1 \geq 1$.

- If S_I is polynomially tractable then $b_d \in \mathcal{O}(\ln d)$ as $d \to \infty$.
- If S_I is strongly polynomially tractable then $b_d \in \mathcal{O}(1)$ as $d \to \infty$, and $\lambda_1 = 1 > \lambda_2$.

Proof. Assume $\lambda_1 \geq 1$ and let τ be given by Lemma 5.6. Then, independently of the amount of symmetry conditions, we have $\lambda_{d,1} = \lambda_1^d \geq 1$ and there exist absolute constants $r \geq 0$ and C > 1 such that

$$\frac{1}{(\lambda_1)^{\tau d}} \sum_{\mathbf{k} \in \nabla_d} (\widetilde{\lambda}_{d,\mathbf{k}})^{\tau} \le C d^r, \quad d \in \mathbb{N},$$
(5.19)

due to Lemma 5.6. In the case of strong polynomial tractability we even have r = 0. For $d \ge 2$ we use the product structure of $\lambda_{d,k}$, $k \in \nabla_d$, provided by (5.14). That is, we split the sum with respect to the coordinates with and without symmetry conditions. Hence,

$$\sum_{\boldsymbol{k}=(\boldsymbol{h},\boldsymbol{j})\in\nabla_{d}}\widetilde{\lambda}_{d,\boldsymbol{k}}^{\tau} = \sum_{\boldsymbol{j}\in(\mathcal{M}_{1})^{b_{d}}}\widetilde{\lambda}_{b_{d},\boldsymbol{j}}^{\tau} \sum_{\substack{\boldsymbol{h}\in(\mathcal{M}_{1})^{a_{d}}\\h_{1}\leq\cdots\leq h_{a_{d}}}}\widetilde{\lambda}_{a_{d},\boldsymbol{h}}^{\tau} = \left(\sum_{m=1}^{\#\mathcal{M}_{1}}\lambda_{m}^{\tau}\right)^{b_{d}} \sum_{\substack{\boldsymbol{h}\in\mathcal{M}_{a_{d}}\\h_{1}\leq\cdots\leq h_{a_{d}}}}\widetilde{\lambda}_{a_{d},\boldsymbol{h}}^{\tau} \quad (5.20)$$

for $d = a_d + b_d \ge 2$, which leads to

$$\left(\sum_{m=1}^{\#\mathcal{M}_1} \left(\frac{\lambda_m}{\lambda_1}\right)^{\tau}\right)^{b_d} \sum_{\substack{\boldsymbol{h} \in \mathcal{M}_{a_d} \\ h_1 \leq \cdots \leq h_{a_d}}} \prod_{l=1}^{a_d} \left(\frac{\lambda_{h_l}}{\lambda_1}\right)^{\tau} \leq C \, d^r.$$

In any case the second sum in the above inequality is bounded from below by 1. Thus, using $\#\mathcal{M}_1 \geq 2$ we conclude that $(1 + \lambda_2^{\tau}/\lambda_1^{\tau})^{b_d} \leq (\sum_{m=1}^{\#\mathcal{M}_1} \lambda_m^{\tau}/\lambda_1^{\tau})^{b_d}$ is polynomially

bounded from above. Since we always assume $\lambda_2 > 0$ this leads to the bounds we claimed for b_d .

It remains to show the assertions on the two largest univariate eigenvalues in the case of strong polynomial tractability. To this end, assume for a moment that $\lambda_1 > 1$. Then, as $\lambda_2 > 0$, there exists some $K \in \mathbb{N}_0$ such that $\lambda_2 \geq (1/\lambda_1)^K$. Now it is easy to see that (independently of the number of symmetry conditions) there are at least $1 + \lfloor d/(K+1) \rfloor$ different $k \in \nabla_d$ such that λ_d is $k \geq 1$. Namely, for $k = 0, \ldots, \lfloor d/(K+1) \rfloor$ we can take the first k = 1 coordinates of $k \in \nabla_d$ equal to one. To the remaining coordinates we assign the value two and obtain

$$\widetilde{\lambda}_{d,\boldsymbol{k}} = \lambda_1^{d-l} \lambda_2^l \ge \lambda_1^{Kl} \lambda_2^l \ge 1.$$

In other words, we have $\lambda_{d,1+\lfloor d/(K+1)\rfloor} \geq 1$. On the other hand, strong polynomial tractability implies $\sum_{i=\lceil 1+C\rceil}^{\infty} \lambda_{d,i}^{\tau} \leq C_1$ for some absolute constants $\tau, C, C_1 > 0$ and all $d \in \mathbb{N}$ (see (5.18)). Consequently, for every $d \geq d_0 = (2+C)(K+1)$ we obtain $1+\lfloor d/(K+1)\rfloor \geq \lceil 1+C\rceil$ and thus

$$C_1 \ge \sum_{i=\lceil 1+C \rceil}^{\infty} \lambda_{d,i}^{\tau} \ge \sum_{i=\lceil 1+C \rceil}^{1+\lfloor d/(K+1) \rfloor} \lambda_{d,i}^{\tau} \ge \lambda_{d,1+\lfloor d/(K+1) \rfloor}^{\tau} (2 + \lfloor d/(K+1) \rfloor - \lceil 1+C \rceil)$$

$$\ge \frac{d}{K+1} - (1+C).$$

Obviously this is a contradiction and we conclude that $\lambda_1 = 1$. Finally, we need to show that necessarily $\lambda_2 < 1$. Assuming $\lambda_1 = \lambda_2 = 1$ leads to K = 0 in the discussion above and hence we obtain the same contradiction as before. Therefore the proof is complete.

Note in passing that independently of the number of symmetry conditions the information complexity $n_{\text{abs}}^{\text{wor}}(\varepsilon, d; S_{d,I_d} \colon \mathcal{B}(\mathfrak{S}_{I_d}(H_d)) \to \mathcal{G}_d)$ grows at least linearly in d if we assume $\lambda_1 \geq 1$ and $\lambda_2 > 0$.

We continue the analysis of I-symmetric problems with respect to the absolute error criterion by proving that the necessary conditions we stated are also sufficient for (strong) polynomial tractability. For this purpose we need a rather technical preliminary lemma. For the convenience of the reader we include a full proof that uses only elementary induction arguments.

LEMMA 5.8. Let $(\mu_m)_{m\in\mathbb{N}}$ be a non-increasing sequence of non-negative real numbers with $\mu_1 > 0$ and set $\mu_{s,\mathbf{k}} = \prod_{l=1}^s \mu_{k_l}$ for $\mathbf{k} \in \mathbb{N}^s$ and $s \in \mathbb{N}$. Then, for all $V \in \mathbb{N}_0$ and every $d \in \mathbb{N}$,

$$\sum_{\substack{\mathbf{k} \in \mathbb{N}^d \\ 1 \le k_1 \le \dots \le k_d}} \mu_{d,\mathbf{k}} \le (\mu_1)^d d^V \left(1 + V + \sum_{L=1}^d (\mu_1)^{-L} \sum_{\substack{\mathbf{j}^{(L)} \in \mathbb{N}^L \\ V + 2 \le j_1^{(L)} \le \dots \le j_L^{(L)}}} \mu_{L,\mathbf{j}^{(L)}} \right). \tag{5.21}$$

Proof. Step 1. By induction on s we first prove that for every fixed $m \in \mathbb{N}$,

$$\sum_{\substack{\boldsymbol{k} \in \mathbb{N}^s \\ m \le k_1 \le \dots \le k_s}} \mu_{s,\boldsymbol{k}} = (\mu_m)^s + \sum_{l=1}^s (\mu_m)^{s-l} \sum_{\substack{\boldsymbol{j}^{(l)} \in \mathbb{N}^l \\ m+1 \le j_1^{(l)} \le \dots \le j_l^{(l)}}} \mu_{l,\boldsymbol{j}^{(l)}} \quad \text{for all } s \in \mathbb{N}.$$
 (5.22)

Easy calculations show that this holds at least for the initial step s = 1. Therefore, assume the assertion (5.22) is true for some $s \in \mathbb{N}$. Then

$$\sum_{\substack{\mathbf{k} \in \mathbb{N}^{s+1} \\ m \le k_1 \le \dots \le k_{s+1}}} \mu_{s+1,\mathbf{k}} = \sum_{k_1 = m}^{\infty} \mu_{k_1} \sum_{\substack{\mathbf{h} \in \mathbb{N}^s \\ k_1 \le h_1 \le \dots \le h_s}} \mu_{s,\mathbf{h}}$$

$$= \mu_m \sum_{\substack{\mathbf{h} \in \mathbb{N}^s \\ m \le h_1 \le \dots \le h_s}} \mu_{s,\mathbf{h}} + \sum_{\substack{\mathbf{k} \in \mathbb{N}^{s+1} \\ m+1 \le k_1 \le \dots \le k_{s+1}}} \mu_{s+1,\mathbf{k}}.$$

Now, by inserting the induction hypothesis for the first sum and renaming k to $j^{(s+1)}$ in the remaining sum, we conclude that $\sum_{k \in \mathbb{N}^{s+1}, m \leq k_1 \leq \cdots \leq k_{s+1}} \mu_{s+1,k}$ equals

$$(\mu_m)^{s+1} + \sum_{l=1}^{s} (\mu_m)^{s+1-l} \sum_{\substack{\boldsymbol{j^{(l)} \in \mathbb{N}^l} \\ m+1 \leq j_1^{(l)} \leq \cdots \leq j_l^{(l)}}} \mu_{l,\boldsymbol{j^{(l)}}} + \sum_{\substack{\boldsymbol{j^{(s+1)} \in \mathbb{N}^{s+1}} \\ m+1 \leq j_1^{(s+1)} \leq \cdots \leq j_{s+1}^{(s+1)}}} \mu_{s+1,\boldsymbol{j^{(s+1)}}}.$$

Hence (5.22) holds for s + 1 as well, and the induction is complete.

Step 2. Here we prove (5.21) via another induction on $V \in \mathbb{N}_0$. Therefore, let $d \in \mathbb{N}$ be arbitrarily fixed. The initial step, V = 0, corresponds to (5.22) for s = d and m = 1. Thus assume (5.21) is valid for some fixed $V \in \mathbb{N}_0$. Then, by using (5.22) for s = L and m = V + 2, we see that the right-hand side of (5.21) equals

$$(\mu_1)^d d^V \Big(1 + V + \sum_{L=1}^d (\mu_1)^{-L} \Big((\mu_{V+2})^L + \sum_{l=1}^L (\mu_{V+2})^{L-l} \sum_{\substack{\boldsymbol{j}^{(l)} \in \mathbb{N}^l \\ (V+2) + 1 \le j_1^{(l)} \le \cdots \le j_l^{(l)}}} \mu_{l,\boldsymbol{j}^{(l)}} \Big) \Big).$$

Now we estimate 1 + V by d(1 + V), take advantage of the non-increasing ordering of $(\mu_m)_{m \in \mathbb{N}}$, and extend the inner sum from L to d in order to obtain

$$\sum_{\substack{\mathbf{k} \in \mathbb{N}^d \\ 1 \le k_1 \le \dots \le k_d}} \mu_{d,\mathbf{k}} \le (\mu_1)^d d^{V+1} \Big(1 + (V+1) + \sum_{l=1}^d (\mu_1)^{-l} \sum_{\substack{\mathbf{j}^{(l)} \in \mathbb{N}^l \\ (V+1) + 2 \le \mathbf{j}_1^{(l)} \le \dots \le \mathbf{j}_l^{(l)}}} \mu_{l,\mathbf{j}^{(l)}} \Big).$$

Since this estimate corresponds to (5.21) for V+1 the claim is proven.

Now the sufficient conditions read as follows.

PROPOSITION 5.9 (Sufficient conditions, symmetric case). Let $P_{I_d} = \mathfrak{S}_{I_d}$, assume S_I is the problem considered in Lemma 5.6, and let $\lambda = (\lambda_m)_{m \in \mathbb{N}} \in \ell_{\tau_0}$ for some $\tau_0 \in (0, \infty)$.

- If $\lambda_1 < 1$ then S_I is strongly polynomially tractable.
- If $\lambda_1 = 1 > \lambda_2$ and $b_d \in \mathcal{O}(1)$ then S_I is strongly polynomially tractable.
- If $\lambda_1 = 1$ and $b_d \in \mathcal{O}(\ln d)$ as $d \to \infty$ then S_I is polynomially tractable.

Proof. Step 1. We start the proof by exploiting the property $\lambda \in \ell_{\tau_0}$; namely we use the ordering of $(\lambda_m)_{m \in \mathbb{N}}$ to conclude that

$$m \lambda_m^{\tau_0} \le \lambda_1^{\tau_0} + \dots + \lambda_m^{\tau_0} < \sum_{i=1}^{\infty} \lambda_i^{\tau_0} = \|\lambda \mid \ell_{\tau_0}\|^{\tau_0} < \infty \quad \text{for any } m \in \mathbb{N}.$$

Hence, there exists some $C_{\tau_0} > 0$ such that λ_m is bounded from above by $C_{\tau_0} m^{-r}$ for every $r \leq 1/\tau_0$. Therefore there is some index such that for every larger $m \in \mathbb{N}$ we have $\lambda_m < 1$. We denote the smallest such index by m_0 . Similar to the calculations of Novak and Woźniakowski [NW08, p. 180] this leads to

$$\sum_{m=m_0}^{\infty} \lambda_m^{\tau} \le (p+1)\lambda_{m_0}^{\tau} + C_{\tau_0}^{\tau} \int_{m_0+p}^{\infty} x^{-\tau r} \, \mathrm{d}\lambda^1(x) = (p+1)\lambda_{m_0}^{\tau} + \frac{C_{\tau_0}^{\tau}}{\tau r - 1} \, \frac{1}{(m_0 + p)^{\tau r - 1}}$$

for every $p \in \mathbb{N}_0$ and all τ such that $\tau r > 1$. In particular, with $r = 1/\tau_0$ we obtain for all $\tau > \tau_0$ and any $p \in \mathbb{N}_0$ the estimate

$$\sum_{m=m_0}^{\infty} (\lambda_m)^{\tau} \le (p+1) (\lambda_{m_0})^{\tau} + \frac{1/\tau}{1/\tau_0 - 1/\tau} \left(\frac{C_{\tau_0}^{1/(1/\tau_0 - 1/\tau)}}{m_0 + p} \right)^{\tau(1/\tau_0 - 1/\tau)}.$$

Note that for a given $\delta > 0$ there exists some constant $\tau_1 \geq \tau_0$ such that for all $\tau > \tau_1$ we have $1/(1/\tau_0 - 1/\tau) \in (\tau_0, \tau_0 + \delta)$. Hence, if $p \in \mathbb{N}_0$ is sufficiently large then we conclude that for all $\tau > \tau_1$,

$$\sum_{m=m_0}^{\infty} (\lambda_m)^{\tau} \le (p+1) (\lambda_{m_0})^{\tau} + \frac{\tau_0 + \delta}{\tau_1} \left(\frac{C_1}{m_0 + p} \right)^{\tau/(\tau_0 + \delta)},$$

where we set $C_1 = \max\{1, C_{\tau_0}^{\tau_0 + \delta}\} < m_0 + p$. Finally, since $\lambda_{m_0} < 1$, both summands tend to zero as $\tau \to \infty$. In particular, there exist $\tau > \tau_1 \ge \tau_0$ such that

$$\sum_{m=m_0}^{\infty} (\lambda_m)^{\tau} \le \frac{1}{2}.$$

Step 2. Now all the assertions stated can be seen using the second point of Theorem 2.8. Indeed, for polynomial tractability it is sufficient to show that

$$\sum_{\mathbf{k} \in \nabla_d} (\widetilde{\lambda}_{d,\mathbf{k}})^{\tau} = \sum_{i=1}^{\infty} (\lambda_{d,i})^{\tau} \le C d^{r\tau} \quad \text{for all } d \in \mathbb{N}$$
 (5.23)

and some $C, \tau > 0$, as well as some $r \ge 0$. If this even holds for r = 0 we obtain strong polynomial tractability.

In the case $\lambda_1 < 1$ we can estimate the sum on the left of (5.23) from above by $(\sum_{m=1}^{\infty} \lambda_m^{\tau})^d$ since clearly $\nabla_d \subseteq \mathcal{M}_d \subseteq \mathbb{N}^d$. Using Step 1 with $m_0 = 1$ we conclude that $\sum_{\mathbf{k} \in \nabla_d} (\widetilde{\lambda}_{d,\mathbf{k}})^{\tau} \leq 2^{-d}$ for some large $\tau > \tau_0$. Hence the problem is strongly polynomially tractable in this case.

For the proof of the remaining points we assume that $\lambda_1 = 1$. In any case we have

$$\sum_{k \in \nabla_1} (\widetilde{\lambda}_{1,k})^{\tau} \le \sum_{m=1}^{\infty} (\lambda_m)^{\tau_0} = \|\lambda \mid \ell_{\tau_0}\|^{\tau_0} < \infty$$

for all $\tau \geq \tau_0$ because $\lambda \in \ell_{\tau_0}$. Therefore we can assume $d \geq 2$ in the following. Recall that we can split the first sum in (5.23) with respect to the coordinates with and without symmetry conditions. That is, for $d = a_d + b_d \geq 2$ we use (5.20).

If $\lambda_2 < 1$ and b_d is universally bounded then the first factor in this splitting can be bounded by a constant and the second factor can be estimated using Lemma 5.8 with

V=0, d replaced by a_d , and μ replaced by λ^{τ} (31). Consequently, for any $\tau \geq \tau_0$,

$$V = 0, \text{ a replaced by } a_d, \text{ and } \mu \text{ replaced by } \lambda^{-}(\gamma). \text{ Consequently, for any } \tau \geq \tau_0,$$

$$\sum_{\substack{\boldsymbol{h} \in \mathcal{M}_{a_d} \\ h_1 \leq \dots \leq h_{a_d}}} (\widetilde{\lambda}_{a_d, \boldsymbol{h}})^{\tau} \leq 1 + \sum_{L=1}^{a_d} \sum_{\substack{\boldsymbol{j}^{(L)} \in \mathbb{N}^L \\ 2 \leq j_1^{(L)} \leq \dots \leq j_L^{(L)}}} (\widetilde{\lambda}_{L, \boldsymbol{j}^{(L)}})^{\tau} \leq 1 + \sum_{L=1}^{a_d} \left(\sum_{m=2}^{\infty} \lambda_m^{\tau}\right)^L. \tag{5.24}$$

Now, with the help of Step 1 and the properties of geometric series, we see that if τ is large enough then (5.24) can be estimated further by $1 + \sum_{L=1}^{\infty} 2^{-L} = 2$. In summary also $\sum_{k \in \nabla_d} (\widetilde{\lambda}_{d,k})^{\tau}$ is universally bounded in this case and therefore the problem S_I is strongly polynomially tractable.

To prove the last point we argue in the same manner. Here the assumption that b_d belongs to $\mathcal{O}(\ln d)$ as $d \to \infty$ implies that the first factor in the splitting (5.20) is polynomially bounded in d. For the second factor we again apply Lemma 5.8, but in this case we set $V = m_0 - 2$, where m_0 denotes the first index $m \in \mathbb{N}$ such that $\lambda_m < 1$. Keep in mind that this index is at least two because of $\lambda_1 = 1$. On the other hand, it is finite, since $\lambda \in \ell_{\tau_0}$. Therefore, by the same arguments as above, the second factor in the splitting (5.20) is polynomially bounded in d, too. All in all, this proves (5.23) and thus S_I is polynomially tractable in this case.

We summarize the results obtained for I-symmetric tensor product problems S_I $(S_{d,I_d})_{d\in\mathbb{N}}$ in the following theorem.

Theorem 5.10 (Tractability of symmetric problems, absolute errors). Let $S_1: H_1 \to \mathcal{G}_1$ denote a compact linear operator between two Hilbert spaces and let $\lambda = (\lambda_m)_{m \in \mathbb{N}}$ be the sequence of eigenvalues of $W_1 = S_1^{\dagger} S_1$ with respect to a non-increasing ordering. Moreover, for d > 1 let $\emptyset \neq I_d \subseteq \{1, \ldots, d\}$ and let $S_I = (S_{d,I_d})_{d \in \mathbb{N}}$ be the linear tensor product problem $S = (S_d)_{d \in \mathbb{N}}$ restricted to the I_d -symmetric subspaces $\mathfrak{S}_{I_d}(H_d)$ of the d-fold tensor product spaces H_d . Consider the worst case setting with respect to the absolute error criterion and let $\lambda_2 > 0$. Then S_I is strongly polynomially tractable if and only if $\lambda \in \ell_{\tau}$ for some $\tau \in (0, \infty)$ and

- $\lambda_1 < 1$, or
- $1 = \lambda_1 > \lambda_2$ and $(d \#I_d) \in \mathcal{O}(1)$ as $d \to \infty$.

Moreover, provided that $\lambda_1 \leq 1$ the problem is polynomially tractable if and only if $\lambda \in \ell_{\tau}$ for some $\tau \in (0, \infty)$ and

- $\lambda_1 < 1$, or
- $\lambda_1 = 1$ and $(d \#I_d) \in \mathcal{O}(\ln d)$ as $d \to \infty$.

Note that we do not have sufficient conditions for polynomial tractability in the case when $\lambda_1 > 1$. We only know that $d - \#I_d \in \mathcal{O}(\ln d)$ as $d \to \infty$ is necessary in this situation. Anyway, we completely characterized strong polynomial tractability of symmetric problems. In this respect we improved the results known from [Wei12a]. Moreover, we have shown that the results hold for finite-dimensional and for non-separable source spaces H_1 as well.

⁽³¹⁾ Observe that this choice implies in particular that $\mu_m = 0$ for any $m > \# \mathcal{M}_1$.

Before we turn to the complexity of antisymmetric problems we briefly focus on the normalized error criterion for the *I*-symmetric setting in the next subsection.

5.3.2. Symmetric problems (normalized errors). Due to (2.10) and (5.14) the information complexity of I-symmetric problems $S_I = (S_{d,I_d})_{d \in \mathbb{N}}$ in the worst case setting with respect to the normalized error criterion is given by

$$n_{\mathrm{norm}}^{\mathrm{wor}}(\varepsilon',d;\mathfrak{S}_{I_d}(H_d)) = \#\bigg\{ \boldsymbol{k} \in \nabla_d \; \bigg| \; \frac{\widetilde{\lambda}_{d,\boldsymbol{k}}}{\lambda_{d,1}} = \prod_{l=1}^d \frac{\lambda_{k_l}}{\lambda_1} > (\varepsilon')^2 \bigg\}$$

for $\varepsilon' \in (0,1)$ and $d \in \mathbb{N}$, since we have $(\varepsilon_d^{\text{init}})^2 = \lambda_{d,1} = \lambda_1^d$ for any kind of symmetric problem. In contrast, for the absolute error, criterion (2.9) yields

$$n_{\mathrm{abs}}^{\mathrm{wor}}(\varepsilon, d; \mathfrak{S}_{I_d}(H_d)) = \# \Big\{ \mathbf{k} \in \nabla_d \mid \widetilde{\lambda}_{d, \mathbf{k}} = \prod_{l=1}^d \lambda_{k_l} > \varepsilon^2 \Big\},$$

where $\varepsilon > 0$ and $d \in \mathbb{N}$. Hence, using the ideas stated in the proof of Theorem 2.12 it suffices to study a scaled tensor product problem $T_d : \mathfrak{S}_{I_d}(H_d) \to \mathcal{G}_d$ with respect to the absolute error criterion in order to obtain tractability results for S_I in the normalized situation. To this end, recall that the squared singular values of T_1 equal $\mu = (\mu_m)_{m \in \mathbb{N}}$ with $\mu_m = \lambda_m/\lambda_1$ such that we always have $\mu_1 = 1$. Furthermore, we obviously have $\mu \in \ell_{\tau}$ if and only if $\lambda \in \ell_{\tau}$. This leads to the following theorem.

THEOREM 5.11 (Tractability of symmetric problems, normalized errors). Consider the situation of Theorem 5.10. We study the worst case setting with respect to the normalized error criterion. Then $S_I = (S_{d,I_d})_{d \in \mathbb{N}}$ is strongly polynomially tractable if and only if

$$\lambda \in \ell_{\tau} \quad \text{for some } \tau \in (0, \infty), \quad \lambda_1 > \lambda_2, \quad d - \#I_d \in \mathcal{O}(1) \quad \text{as } d \to \infty.$$

Moreover, the problem S_I is polynomially tractable if and only if

$$\lambda \in \ell_{\tau}$$
 for some $\tau \in (0, \infty)$ and $d - \#I_d \in \mathcal{O}(\ln d)$ as $d \to \infty$.

5.3.3. Antisymmetric problems (absolute errors). We start with sufficient conditions for (strong) polynomial tractability which slightly improve the results stated in [Wei12a, Proposition 5].

PROPOSITION 5.12 (Sufficient conditions, antisymmetric case). Let $P_{I_d} = \mathfrak{A}_{I_d}$, let $S_I = (S_{d,I_d})_{d \in \mathbb{N}}$ be the problem considered in Lemma 5.6, and let $\lambda = (\lambda_m)_{m \in \mathbb{N}} \in \ell_{\tau_0}$ for some $\tau_0 \in (0,\infty)$.

- If $\lambda_1 < 1$ then S_I is strongly polynomially tractable, independent of the number of antisymmetry conditions.
- If $\lambda_1 \geq 1$ and if there exist constants $\tau \geq \tau_0$, $d_0 \in \mathbb{N}$, as well as $C \geq 1$, and $q \geq 0$ such that the number of antisymmetric coordinates a_d in dimension d satisfies

$$\frac{\ln(a_d!)}{d} + \frac{\ln(Cd^q)}{d} \ge \ln(\|\lambda \mid \ell_\tau\|^\tau) \quad \text{for all } d \ge d_0$$
 (5.25)

then the problem S_I is polynomially tractable. If this even holds for q=0 then we obtain strong polynomial tractability.

Proof. Just as for the symmetric setting, the proof of these sufficient conditions is based on the second point of Theorem 2.8. We show that under the given assumptions for some $\tau \geq \tau_0$ the whole sum of the eigenvalues

$$\sum_{\mathbf{k} \in \nabla_d} (\widetilde{\lambda}_{d,\mathbf{k}})^{\tau} = \sum_{i=1}^{\infty} (\lambda_{d,i})^{\tau}$$
(5.26)

is universally bounded, or polynomially bounded in d, respectively. Note that since we deal with the case $P_{I_d} = \mathfrak{A}_{I_d}$ now, the set ∇_d is given by the second line in (5.9). Moreover observe that for d=1 there is no antisymmetry condition at all. That is, we have $\nabla_1 = \mathcal{M}_1 \subseteq \mathbb{N}$ and the sums in (5.26) equal $\|\lambda \mid \ell_\tau\|^\tau \leq \|\lambda \mid \ell_{\tau_0}\|^\tau$ in this case. Therefore, due to the hypothesis $\lambda \in \ell_{\tau_0}$, the term for d=1 is finite.

Hence, let $d \geq 2$ be arbitrarily fixed. Without loss of generality we may reorder the set of coordinates so that $I_d = \{i_1, \ldots, i_{a_d}\} = \{1, \ldots, a_d\}$. That means we assume partial antisymmetry with respect to the first a_d coordinates. For $s \in \mathbb{N}$ with $s \geq d$ let us define cubes of multi-indices

$$Q_{d,s} = \{1, \dots, s\}^d.$$

Furthermore, let $U_{a_d,s} = \{ j \in Q_{a_d,s} \mid j_1 < \dots < j_{a_d} \}$ denote the a_d -dimensional projection of $Q_{d,s}$ which reflects the antisymmetry conditions we assume. With this notation we obtain

$$\sum_{\boldsymbol{k} \in \nabla_d} (\widetilde{\lambda}_{d,\boldsymbol{k}})^\tau = \lim_{s \to \infty} \sum_{\boldsymbol{k} \in \nabla_d \cap Q_{d,s}} (\widetilde{\lambda}_{d,\boldsymbol{k}})^\tau,$$

where the set of multi-indices under consideration $\nabla_d \cap Q_{d,s}$ can be represented as a subset of $U_{a_d,s} \times Q_{b_d,s}$. We will assume $b_d = d - a_d > 0$ in what follows to ensure that this splitting is non-trivial. By the product structure of $\widetilde{\lambda}_{d,k}$, $k \in \nabla_d$, this implies

$$\sum_{\mathbf{k}=(\mathbf{j},\mathbf{i})\in\nabla_{d}\cap Q_{d,s}} (\widetilde{\lambda}_{d,\mathbf{k}})^{\tau} \leq \left(\sum_{\mathbf{j}\in U_{a_{d},s}} \prod_{l=1}^{a_{d}} \lambda_{j_{l}}^{\tau}\right) \left(\sum_{\mathbf{i}\in Q_{b_{d},s}} \prod_{l=1}^{b_{d}} \lambda_{i_{l}}^{\tau}\right).$$
(5.27)

Since the sequence $\lambda = (\lambda_m)_{m \in \mathbb{N}}$ is an element of $\ell_{\tau_0} \hookrightarrow \ell_{\tau}$, we can easily estimate the second factor for every $s \geq d$ from above by

$$\sum_{i \in Q_{b,i,s}} \prod_{l=1}^{b_d} \lambda_{i_l}^{\tau} = \prod_{l=1}^{b_d} \sum_{m=1}^{s} \lambda_m^{\tau} = \left(\sum_{m=1}^{s} \lambda_m^{\tau}\right)^{b_d} \le \left(\sum_{m=1}^{\infty} \lambda_m^{\tau}\right)^{(1/\tau)b_d \tau} = \|\lambda \| \ell_{\tau}\|^{b_d \tau}.$$
 (5.28)

To handle the first term we need an additional argument. Note that due to the structure of $U_{a_d,s}$ we have

$$\sum_{\substack{\boldsymbol{j} \in Q_{a_d,s} \\ \exists k,m: \, j_k = j_m}} \prod_{l=1}^{a_d} \lambda_{j_l}^{\tau} = \sum_{\substack{\boldsymbol{j} \in Q_{a_d,s} \\ \exists k,m: \, j_k = j_m}} \prod_{l=1}^{a_d} \lambda_{j_l}^{\tau} + a_d! \sum_{\substack{\boldsymbol{j} \in U_{a_d,s} \\ \mid l=1}} \prod_{l=1}^{a_d} \lambda_{j_l}^{\tau} \geq a_d! \sum_{\substack{\boldsymbol{j} \in U_{a_d,s} \\ \mid l=1}} \prod_{l=1}^{a_d} \lambda_{j_l}^{\tau}.$$

Consequently, using the same arguments as in (5.28), this yields the upper bound

$$\|\lambda \mid \ell_{\tau}\|^{a_d \tau}/(a_d!)$$

for the first factor in (5.27). Once again this bound does not depend on $s \geq d$. Hence,

due to $d = a_d + b_d$, we conclude that

$$\sum_{\boldsymbol{k} \in \nabla_d} (\widetilde{\lambda}_{d,\boldsymbol{k}})^{\tau} = \lim_{s \to \infty} \sum_{\boldsymbol{k} \in \nabla_d \cap Q_{d,s}} (\widetilde{\lambda}_{d,\boldsymbol{k}})^{\tau} \le \frac{1}{a_d!} \|\lambda \mid \ell_{\tau}\|^{\tau d} \quad \text{ for every } d \in \mathbb{N}$$

and any choice of \mathfrak{A}_{I_d} . Of course, for every $d < d_0$ this upper bound is trivially less than an absolute constant. Thus, to prove the second assertion of Proposition 5.12 it is enough to show that

$$\frac{1}{a_d!} \|\lambda \mid \ell_\tau\|^{\tau d} \le C d^q \quad \text{ for all } d \ge d_0,$$

as well as for some $C \geq 1$ and some $q \geq 0$. But this is equivalent to our hypothesis stated in (5.25). Hence the condition (5.25) implies (strong) polynomial tractability of S_I , independently of the value of λ_1 .

Note that now it suffices to show that $\lambda_1 < 1$ already yields (5.25) with q = 0 and C = 1 in order to complete the proof. To see this, observe that (due to Step 1 in the proof of Proposition 5.9) we know that there exists some $\tau > \tau_0$ such that $\|\lambda\| \ell_{\tau}\|^{\tau} = \sum_{m=1}^{\infty} \lambda_m^{\tau}$ is strictly less than 1. Thus the right-hand side of (5.25) is negative in this case, whereas the left-hand side is non-negative for every choice of a_d .

Let us briefly comment on the above result. Clearly, for any $q \geq 0$ the term $\ln(C d^q)/d$ in (5.25) tends to zero as $d \to \infty$. Hence there is not much difference in the sufficient conditions we stated for strong polynomial and for polynomial tractability. Moreover, we mention that Theorem 2.8 allows us to omit the f(d) - 1 largest eigenvalues $\lambda_{d,i}$, where f(d) may grow polynomially in $(\varepsilon_d^{\text{init}})^{-1}$ with d, but we did not use this fact in the above proof.

The next example investigates how fast a_d needs to grow with the dimension d in order to fulfill condition (5.25).

EXAMPLE 5.13. For any $d \in \mathbb{N}$ and some $\gamma > 0$ let

$$a_d = \left\lceil \frac{d}{\ln d^{\gamma}} \right\rceil. \tag{5.29}$$

Then Stirling's formula provides $a_d \ln(a_d/e) \le \ln(a_d!) < \ln a_d - a_d + a_d \ln a_d$ if d (and hence also a_d) is sufficiently large. Consequently,

$$\frac{\ln(a_d!)}{d} \ge \frac{a_d \ln(a_d/e)}{d} \ge \frac{1}{\gamma} \frac{\ln\left(d \cdot \frac{1}{e\gamma \ln d}\right)}{\ln d} = \frac{1}{\gamma} \left(1 - \frac{\ln(e\gamma \ln d)}{\ln d}\right) \nearrow \frac{1}{\gamma}$$

as $d \to \infty$. On the other hand, we have $a_d/d \le 1/(\gamma \ln d) + 1/d$ and thus

$$\frac{\ln(a_d!)}{d} < \frac{\ln a_d - a_d}{d} + \left(\frac{1}{\gamma \ln d} + \frac{1}{d}\right) \ln a_d = \frac{2 \ln a_d - a_d}{d} + \frac{1}{\gamma} \frac{\ln a_d}{\ln d} \le \frac{1}{\gamma}.$$

So we see that γ in (5.29) needs to be strictly smaller than $\ln^{-1}(\|\lambda \mid \ell_{\tau}\|^{\tau})$ in order to fulfill (5.25) with q = 0. In particular, it follows that assumptions like $a_d = \lceil d^{\beta} \rceil$ with $\beta < 1$ are not sufficient to imply tractability using the second point of Proposition 5.12. \square

Now we turn to necessary conditions. As in the symmetric setting Lemma 5.6 shows that $\lambda \in \ell_{\tau}$ is needed for polynomial tractability. In addition, we will see that we need a condition similar to (5.25), particularly if we deal with slowly decreasing eigenvalues λ .

PROPOSITION 5.14 (Necessary conditions, antisymmetric case). Let $P_{I_d} = \mathfrak{A}_{I_d}$ and assume $S_I = (S_{d,I_d})_{d \in \mathbb{N}}$ is the problem considered in Lemma 5.6. Furthermore, let S_I be polynomially tractable with constants C, p > 0 and $q \geq 0$. Then, as $d \to \infty$, the initial error $\varepsilon_d^{\text{init}}$ tends to zero faster than the inverse of any polynomial. Moreover, $\lambda = (\lambda_m)_{m \in \mathbb{N}} \in \ell_\tau$ for every $\tau > p/2$ and there exist $d^* \in \mathbb{N}$ and $C_2 \geq 1$ such that

$$\frac{1}{d} \sum_{m=1}^{a_d} \ln \frac{\|\lambda \mid \ell_\tau\|^\tau}{\lambda_m^\tau} + \frac{\ln(C_2 d^{2q\tau/p})}{d} \ge \ln(\|\lambda \mid \ell_\tau\|^\tau) \quad \text{for all } d \ge d^*.$$
 (5.30)

Thus we have either $\lambda_1 < 1$ or $\lim_{d \to \infty} a_d = \infty$.

Proof. Step 1. For the whole proof assume $\tau > p/2$ is fixed. Then Lemma 5.6 shows that $\lambda \in \ell_{\tau}$. As in (5.20) for the symmetric case, we can split the sum of the eigenvalues so that for all $d \in \mathbb{N}$,

$$\sum_{\boldsymbol{k} \in \nabla_d} (\widetilde{\lambda}_{d,\boldsymbol{k}})^{\tau} = \left(\sum_{m=1}^{\#\mathcal{M}_1} \lambda_m^{\tau}\right)^{b_d} \sum_{\substack{\boldsymbol{j} \in \mathcal{M}_{a_d} \\ j_1 < \dots < j_{a_d}}} (\widetilde{\lambda}_{a_d,\boldsymbol{j}})^{\tau} \ge \|\lambda \| \ell_{\tau}\|^{\tau b_d} \cdot \lambda_1^{\tau} \cdot \dots \cdot \lambda_{a_d}^{\tau}.$$

Hence Lemma 5.6 together with the fact that $\lambda_{d,1} = \lambda_1^{b_d} \cdot \lambda_1 \cdot \ldots \cdot \lambda_{a_d}$ gives

$$\left(\frac{\|\lambda \mid \ell_{\tau}\|^{\tau}}{\lambda_{1}^{\tau}}\right)^{b_{d}} \leq (1+C) d^{q} + C^{2\tau/p} \zeta\left(\frac{2\tau}{p}\right) \left(\frac{d^{2q/p}}{\lambda_{d,1}}\right)^{\tau}.$$
(5.31)

We will use this inequality to deduce all the assertions stated in Proposition 5.14.

Step 2. Here we prove the limit property for the initial error $\varepsilon_d^{\text{init}} = \sqrt{\lambda_{d,1}}$, i.e. we show that for every fixed polynomial $\mathcal{P} > 0$,

$$\lambda_{d,1} \cdot \mathcal{P}(d) \to 0 \quad \text{as } d \to \infty.$$
 (5.32)

Since $\lambda_{d,1} \leq \lambda_1^{b_d} \lambda_1^{a_d} = \lambda_1^d$ we can restrict ourselves to the non-trivial case $\lambda_1 \geq 1$ in the following. Assume that there exists a subsequence $(d_k)_{k \in \mathbb{N}}$ of natural numbers, as well as some constant $C_0 > 0$, such that $\lambda_{d_k,1} \mathcal{P}(d_k)$ is bounded from below by C_0 for every $k \in \mathbb{N}$. Then for every $d = d_k$ the right-hand side of (5.31) is bounded from above by some other polynomial $\mathcal{P}_1(d_k) > 0$. On the other hand, due to the general condition $\lambda_2 > 0$, the term $\|\lambda \mid \ell_\tau\|^\tau / \lambda_1^\tau$ is strictly larger than one. Thus it follows that there exists some $C_1 > 0$ such that

$$b_{d_k} \le C_1 \ln d_k$$
 for every $k \in \mathbb{N}$.

Therefore $a_{d_k}=d_k-b_{d_k}\to\infty$ as $k\to\infty$. Moreover, the assumed boundedness of $\lambda_{d_k,1}\mathcal{P}(d_k)$ leads to

$$C_0 \mathcal{P}(d_k)^{-1} \le \lambda_{d_k,1} \le \lambda_1^{C_1 \ln d_k} \cdot \lambda_1 \cdot \ldots \cdot \lambda_{a_{d_k}} = d_k^{C_1 \ln \lambda_1} \cdot \lambda_1 \cdot \ldots \cdot \lambda_{a_{d_k}}$$

since $\lambda_1 \geq 1$. In the first step of the proof of Proposition 5.9 we saw that $\lambda \in \ell_{\tau}$ yields the existence of some $C_{\tau} > 0$ such that $\lambda_m \leq C_{\tau} m^{-1/\tau}$ for every $m \in \mathbb{N}$. Indeed, this

holds for $C_{\tau} = \|\lambda \mid \ell_{\tau}\| > 1$. Hence $\lambda_1^{\tau} \cdot \ldots \cdot \lambda_{a_{d_k}}^{\tau} \leq C_{\tau}^{\tau a_{d_k}} (a_{d_k}!)^{-1}$, which gives $(a_{d_k}/e)^{a_{d_k}} \leq a_{d_k}! \leq (C_{\tau}^{\tau})^{a_{d_k}} \mathcal{P}_2(d_k)$ for all $k \in \mathbb{N}$

and some other polynomial $\mathcal{P}_2 > 0$. If k is sufficiently large then we conclude that

$$a_{d_k} \le a_{d_k} \ln \frac{a_{d_k}}{e C_{\tau}^{\tau}} \le \ln \mathcal{P}_2(d_k),$$

since $a_{d_k} \to \infty$ implies $a_{d_k}/(e C_{\tau}^{\tau}) \ge e$ for $k \ge k_0$. Therefore the number of antisymmetric coordinates a_d needs to be logarithmically bounded from above for every d from the sequence $(d_k)_{k \ge k_0}$. Because b_{d_k} was also found to be logarithmically bounded this contradicts $d_k = a_{d_k} + b_{d_k}$. Consequently, the hypothesis $\lambda_{d_k,1} \mathcal{P}(d_k) \ge C_0 > 0$ cannot be true for any subsequence $(d_k)_k$. In other words, (5.32) holds.

Step 3. Next we show (5.30). From Step 2 we know that there exists $d^* \in \mathbb{N}$ such that $1/\lambda_{d,1} \geq 1$ for all $d \geq d^*$. Hence, (5.31) together with $\tau > p/2$ implies

$$\left(\frac{\|\lambda \mid \ell_{\tau}\|^{\tau}}{\lambda_{1}^{\tau}}\right)^{b_{d}} \leq C_{2} \left(\frac{d^{2q/p}}{\lambda_{d,1}}\right)^{\tau} = \frac{C_{2} d^{2q\tau/p}}{\lambda_{1}^{\tau b_{d}} \cdot \lambda_{1}^{\tau} \cdot \dots \cdot \lambda_{a_{d}}^{\tau}} \quad \text{for } d \geq d^{*},$$

where we set $C_2 = 1 + C + C^{2\tau/p} \zeta(2\tau/p)$. Therefore we obtain

$$C_2 d^{2q\tau/p} \prod_{k=1}^{a_d} \frac{\|\lambda \mid \ell_\tau\|^\tau}{\lambda_k^\tau} \ge \|\lambda \mid \ell_\tau\|^{\tau d}$$

for all $d \ge d^*$, which is equivalent to the estimate (5.30) we claimed.

Step 4. It remains to show that $\lambda_1 \geq 1$ implies that $\lim_{d \to \infty} a_d$ is infinite. To this end, note that every summand in (5.30) is strictly positive. If we assume for a moment the existence of a subsequence $(d_k)_{k \in \mathbb{N}}$ such that a_{d_k} is bounded for every $k \in \mathbb{N}$ then the left-hand side of (5.30) is less than some positive constant divided by d_k . Hence it tends to zero as $k \to \infty$. On the other hand, the right-hand side of (5.30) is strictly larger than some positive constant, since $\lambda_1 \geq 1$ and $\lambda_2 > 0$. This contradiction completes the proof.

As mentioned before, there are examples such that the sufficient condition (5.25) from Proposition 5.12 is also necessary (up to some constant factor) in order to conclude polynomial tractability in the antisymmetric setting. Now we are ready to give such an example.

EXAMPLE 5.15. Consider the situation of Lemma 5.6 for $P_{I_d} = \mathfrak{A}_{I_d}$ and assume the problem S_I is polynomially tractable. In addition, for a fixed $\tau \in (0, \infty)$, let $\lambda = (\lambda_m)_{m \in \mathbb{N}} \in \ell_{\tau}$ with $\lambda_1 \geq 1$ and assume the existence of some $m_0 \in \mathbb{N}$ such that

$$\lambda_m \ge \frac{\|\lambda \mid \ell_\tau\|}{m^{\alpha/\tau}}$$
 for all $m > m_0$ and some $\alpha > 1$. (5.33)

Then we claim that there exist constants $\bar{d} \in \mathbb{N}$, $C \geq 1$, and $r \geq 0$ such that

$$\alpha \frac{\ln(a_d!)}{d} + \frac{\ln(Cd^r)}{d} \ge \ln(\|\lambda \mid \ell_\tau\|^\tau) \quad \text{for all } d \ge \bar{d}.$$
 (5.34)

Recall that due to Proposition 5.12, for the degree of antisymmetry a_d , it was sufficient to assume (5.34) with $\alpha = 1$ in order to conclude (strong) polynomial tractability (see (5.25)). Moreover, keep in mind that we know from Example 5.13 that $\ln(a_d!)/d$ tends to $1/\gamma$ if we assume that a_d is given by (5.29). Hence in the present example we have strong polynomial tractability if $\gamma < \ln^{-1}(\|\lambda + \ell_{\tau}\|^{\tau})$, whereas the problem is polynomially intractable if $\gamma > \alpha/\ln(\|\lambda + \ell_{\tau}\|^{\tau})$.

To verify (5.34) we can use Proposition 5.14 and, in particular, inequality (5.30). Since $\lambda_1 \geq 1$ we know that $\lim_d a_d = \infty$, i.e. $a_d > m_0$ for every d larger than some $d_1 \in \mathbb{N}$. Furthermore, note that (5.33) is equivalent to

$$\ln \frac{\|\lambda \mid \ell_{\tau}\|^{\tau}}{\lambda_{m}^{\tau}} \leq \alpha \ln m \quad \text{ for all } m > m_{0}.$$

Hence if $d \ge d_1$ then we can estimate the sum in (5.30) from above by

$$\frac{1}{d} \sum_{m=1}^{a_d} \ln \frac{\|\lambda \mid \ell_{\tau}\|^{\tau}}{\lambda_m^{\tau}} \leq \frac{m_0}{d} \ln \frac{\|\lambda \mid \ell_{\tau}\|^{\tau}}{\lambda_{m_0}^{\tau}} + \frac{\alpha}{d} \sum_{m=m_0+1}^{a_d} \ln m \leq \frac{C_{\lambda}}{d} + \alpha \frac{\ln(a_d!)}{d}.$$

Obviously, for d larger than some $d_2 \in \mathbb{N}$ the term $C_{\lambda} + \ln(C_2 d^{2q\tau/p})$ is less than $\ln(C d^r)$, where $C \geq 1$ and $r \geq 0$. Here r = 0 if and only if q = 0 in (5.14), i.e. if the problem is strongly polynomially tractable. Consequently, we can derive (5.34) from (5.30) by choosing $\bar{d} = \max\{d_1, d_2, d^*\}$. \square

Although there remains a small gap between the necessary and the sufficient conditions for the absolute error criterion, the most important cases of antisymmetric tensor product problems are covered by our results. Let us summarize the main facts.

THEOREM 5.16 (Tractability of antisymmetric problems, absolute errors). Let $S_1: H_1 \to \mathcal{G}_1$ denote a compact linear operator between two Hilbert spaces and let $\lambda = (\lambda_m)_{m \in \mathbb{N}}$ be the sequence of eigenvalues of $W_1 = S_1^{\dagger} S_1$ with non-increasing ordering. Moreover, for d > 1 let $\emptyset \neq I_d \subseteq \{1, \ldots, d\}$ and assume $S_I = (S_{d,I_d})_{d \in \mathbb{N}}$ is the linear tensor product problem $S = (S_d)_{d \in \mathbb{N}}$ restricted to the I_d -antisymmetric subspaces $\mathfrak{A}_{I_d}(H_d)$ of the d-fold tensor product spaces H_d . Consider the worst case setting with respect to the absolute error criterion and let $\lambda_2 > 0$. Then for the case $\lambda_1 < 1$ the following statements are equivalent:

- S_I is strongly polynomially tractable.
- S_I is polynomially tractable.
- There exists a constant $\tau \in (0, \infty)$ such that $\lambda \in \ell_{\tau}$.

Moreover, the same equivalences hold true if $\lambda_1 \geq 1$ and $\#I_d$ grows linearly with d.

At this point we mention that for fully antisymmetric problems, i.e. $\#I_d = a_d = d$, an explicit formula for the information complexity with respect to the absolute error criterion is known. Furthermore, simple examples can be constructed which show that

we cannot expect the same nice tractability behavior if we deal with normalized errors. For further details the interested reader is referred to [Wei11, Proposition 8].

5.3.4. Antisymmetric problems (normalized errors). Up to now every complexity assertion in this chapter was mainly based on Theorem 2.8, which dealt with the general situation of arbitrary compact linear operators between Hilbert spaces and with the absolute error criterion. While investigating tractability properties of I-symmetric problems with respect to the normalized error criterion, we were able to use assertions from the absolute error setting. Since for I-antisymmetric problems the structure of the initial error is more complicated, this approach will not work again. Therefore we recall Theorem 2.9 as a replacement of Theorem 2.8 for the normalized setting. Based on this we can give the following necessary conditions for (strong) polynomial tractability.

PROPOSITION 5.17 (Necessary conditions, antisymmetric case). Let $S_I = (S_{d,I_d})_{d \in \mathbb{N}}$ denote an I-antisymmetric problem as defined at the beginning of Section 5.3 and consider the worst case setting with respect to normalized errors. Then the fact that S_I is polynomially tractable with constants C, p > 0 and $q \ge 0$ implies that $\lambda = (\lambda_m)_{m \in \mathbb{N}} \in \ell_{\tau}$ for all $\tau > p/2$. Moreover, as $d \to \infty$, $\varepsilon_d^{\text{init}}$ tends to zero faster than the inverse of any polynomial and $b_d \in \mathcal{O}(\ln d)$ as $d \to \infty$. Thus $\lim_{d \to \infty} a_d/d = 1$. In addition, if S_I is strongly polynomially tractable then $b_d \in \mathcal{O}(1)$ as $d \to \infty$.

Proof. From Theorem 2.9 it follows that there is some $C_1 > 0$ such that

$$\frac{1}{(\lambda_{d,1})^{\tau}} \sum_{\mathbf{k} \in \nabla_d} (\widetilde{\lambda}_{d,\mathbf{k}})^{\tau} = \sum_{i=1}^{\infty} \left(\frac{\lambda_{d,i}}{\lambda_{d,1}} \right)^{\tau} \le C_1 d^{2\tau q/p} \quad \text{for every } d \in \mathbb{N}$$
 (5.35)

and all $\tau > p/2$. Once more the index set ∇_d is given as in (5.9). Indeed, Theorem 2.9 shows that it is sufficient to take $C_1 = 2 (1+C)^{2\tau/p} \zeta(2\tau/p)$. As in the proof of Lemma 5.6 it suffices to consider the case d = 1 in (5.35) to see that $\lambda \in \ell_{\tau}$ is necessary for polynomial tractability. Moreover, as in Step 1 of the proof of Proposition 5.14, it follows that

$$\left(\frac{\|\lambda \mid \ell_{\tau}\|^{\tau}}{\lambda_{1}^{\tau}}\right)^{b_{d}} \le C_{1} d^{2\tau q/p}, \quad d \in \mathbb{N},$$
(5.36)

since $\lambda_{d,1} = \lambda_1^{b_d} \cdot \lambda_1 \cdot \ldots \cdot \lambda_{a_d}$. Due to the general assertion $\lambda_2 > 0$ we have $\|\lambda \mid \ell_\tau\|^\tau > \lambda_1^\tau$, and thus polynomial tractability of S_I implies $b_d \leq C_2 \ln d$ for some $C_2 \geq 0$, i.e. b_d is $\mathcal{O}(\ln d)$ as $d \to \infty$. Therefore we obviously have

$$1 \ge \frac{a_d}{d} = 1 - \frac{b_d}{\ln d} \frac{\ln d}{d} \ge 1 - C_2 \frac{\ln d}{d} \to 1, \quad d \to \infty.$$

The proof that strong polynomial tractability leads to $b_d \in \mathcal{O}(1)$ as $d \to \infty$ can be obtained using (5.36) with the same arguments as before and q = 0. Finally we need to show the assertion concerning $\varepsilon_d^{\text{init}}$. Here we refer to Step 2 in the proof of Proposition 5.14.

5.4. Applications. This last section is devoted to applications of the theory developed previously. In Section 5.4.1 we follow the introduction of [Wei12a] and illustrate the power of additional (anti)symmetry conditions on linear tensor product problems by using simple toy examples. Afterwards, in Section 5.4.2, we direct our attention to more advanced problems which we are faced with in practice. There we briefly introduce wavefunctions

and show how our results allow one to handle the approximation problem for such classes of functions.

5.4.1. Toy examples. The aim of the following simple examples is to show that exploiting an a priori knowledge about (anti)symmetries of a given tensor product problem can help to obtain tractability, but it does not make the problem trivial in general.

Let $S = (S_d : H_d \to \mathcal{G}_d)_{d \in \mathbb{N}}$ denote a tensor product problem between Hilbert spaces. Remember that by Section 2.4 for complexity studies it suffices to specify the singular values of the univariate operator S_1 . To simplify the presentation we slightly abuse the notation and denote the information complexity of the entire problem S by $n^{\text{ent}}(\varepsilon, d)$. We want to compare this quantity with the respective information complexities of the restrictions of S to the fully symmetric and the fully antisymmetric subspaces of $(H_d)_{d \in \mathbb{N}}$. These numbers will be denoted by $n^{\text{sym}}(\varepsilon, d)$ and $n^{\text{asy}}(\varepsilon, d)$, respectively.

Clearly, our results show that in any case (as long as we deal with the worst case setting and the absolute error criterion)

$$n^{\mathrm{asy}}(\varepsilon, d) \leq n^{\mathrm{sym}}(\varepsilon, d) \leq n^{\mathrm{ent}}(\varepsilon, d)$$
 for every $\varepsilon > 0$ and all $d \in \mathbb{N}$,

where for d = 1 the terms coincide, since then we do not claim any (anti)symmetry. To see that additional (anti)symmetry conditions may reduce the information complexity dramatically, consider the following three examples.

EXAMPLE 5.18. Let us have a look at the simple case of a linear operator S_1 with singular values σ such that $\lambda_1 = \lambda_2 = 1$ and $\lambda_j = 0$ for $j \geq 3$. Then the information complexity of the entire tensor product problem can be shown to be

$$n^{\text{ent}}(\varepsilon, d) = 2^d$$
 for all $d \in \mathbb{N}$ and $\varepsilon < 1$.

Hence the problem suffers from the curse of dimensionality and is therefore intractable.

On the other hand, our results show that in the fully symmetric setting we have polynomial tractability, because

$$n^{\text{sym}}(\varepsilon, d) = d + 1$$
 for all $d \in \mathbb{N}$ and $\varepsilon < 1$.

Moreover, it can be proved that in this case the complexity of the fully antisymmetric problem decreases with increasing dimension d and finally the problem even gets trivial. In detail, we have

$$n^{\mathrm{asy}}(\varepsilon,d) = \max\left\{3-d,0\right\} \quad \text{ for all } d \in \mathbb{N} \text{ and } \varepsilon < 1,$$

which yields strong polynomial tractability.

□

EXAMPLE 5.19. Next let us consider a more challenging problem, where $\lambda_1 = \cdots = \lambda_m = 1$ and $\lambda_j = 0$ for every $j > m \ge 2$. For m = 2 this obviously coincides with the example studied above, but letting m increase may tell us more about the structure of (anti)symmetric tensor product problems. In this situation it is easy to check that for every $d \in \mathbb{N}$ and all $\varepsilon < 1$,

$$n^{\mathrm{ent}}(\varepsilon, d) = m^d$$
 and $n^{\mathrm{asy}}(\varepsilon, d) = \begin{cases} \binom{m}{d}, & d \leq m, \\ 0, & d > m. \end{cases}$

Since $\binom{m}{d} \geq 2^{d-1}$ for $d \leq \lfloor m/2 \rfloor$, this means that for large m the complexity in the antisymmetric case increases exponentially fast with d up to a certain maximum. Beyond this point it falls back to zero.

The information complexity in the symmetric setting is much harder to calculate for this case. However, it can be seen that we have polynomial tractability, but $n^{\text{sym}}(\varepsilon,d)$ needs to grow at least linearly with d so that the symmetric problem cannot be strongly polynomially tractable, whereas this holds in the antisymmetric setting. The entire problem again suffers from the curse of dimensionality. \Box

EXAMPLE 5.20. For a last illustrative example consider the case $\lambda_1 = 1$ and $\lambda_{j+1} = j^{-\beta}$ for some $\beta \geq 0$ and all $j \in \mathbb{N}$. That means, we have the two largest singular values $\sigma_1 = \sigma_2$ of S_1 equal to one. The remaining series decays like the inverse of some polynomial. If $\beta = 0$ then the operator S_1 is not compact, since the sequence $\lambda = (\lambda_m)_{m \in \mathbb{N}}$ does not tend to zero; hence all the information complexities are infinite in this case.

For $\beta > 0$, any $\delta > 0$, and some C > 0 we have

$$n^{\mathrm{ent}}(\varepsilon, d) \ge 2^d$$
, $n^{\mathrm{sym}}(\varepsilon, d) \ge d + 1$, and $n^{\mathrm{asy}}(\varepsilon, d) \le C\varepsilon^{-(2/\beta + \delta)}$,

for all $\varepsilon < 1$ as well as every $d \in \mathbb{N}$. Thus, again for the entire problem we observe the curse, whereas the antisymmetric problem is strongly polynomially tractable. Once more, the symmetric problem can be shown to be polynomially tractable. Note that in this example the antisymmetric case is not trivial, because all λ_j are strictly positive. If we replace $j^{-\beta}$ by $\log^{-1}(j+1)$ in this example we obtain (polynomial) intractability even in the antisymmetric setting. \square

5.4.2. Wavefunctions. During the last few decades there has been considerable interest in finding approximations of so-called *wavefunctions*, e.g., solutions of the electronic Schrödinger equation. Due to the *Pauli principle* of quantum physics only functions with certain (anti)symmetry properties are of physical interest. For a more detailed view see, e.g., Hamaekers [Ham09], Yserentant [Yse10], or Zeiser [Zei10]. Furthermore, for a comprehensive introduction to the topic, as well as a historical survey, we refer the reader to Hunziker and Sigal [HS00] and Reed and Simon [RS78].

In particular, the notion of multiple partial antisymmetry with respect to two sets of coordinates is useful for describing wavefunctions Ψ . In computational chemistry such functions occur as models which describe quantum states of certain physical d-particle systems. Formally, these functions depend on d blocks of variables $\mathbf{y}_i = (\mathbf{x}^{(i)}, s^{(i)})$, for $i = 1, \ldots, d$, which represent the spacial coordinates $\mathbf{x}^{(i)} = (x_1^{(i)}, x_2^{(i)}, x_3^{(i)}) \in \mathbb{R}^3$ and certain additional intrinsic parameters $s^{(i)} \in C$ of each particle \mathbf{y}_i within the system. Hence, rearranging the arguments so that $\mathbf{x} = (\mathbf{x}^{(1)}, \ldots, \mathbf{x}^{(d)})$ and $\mathbf{s} = (s^{(1)}, \ldots, s^{(d)})$ yields

$$\Psi \colon (\mathbb{R}^3)^d \times C^d \to \mathbb{R}, \quad (\boldsymbol{x}, \boldsymbol{s}) \mapsto \Psi(\boldsymbol{x}, \boldsymbol{s}).$$

In the case of systems of electrons one of the most important parameters is called *spin* and it can take only two values, i.e., $s^{(i)} \in C = \{-1/2, +1/2\}$. Due to the Pauli principle the only wavefunctions Ψ that are physically admissible are those which are antisymmetric

in the sense that for $I \subseteq \{1, \ldots, d\}$ and $I^c = \{1, \ldots, d\} \setminus I$,

$$\Psi(\pi(\boldsymbol{x}), \pi(\boldsymbol{s})) = (-1)^{|\pi|} \Psi(\boldsymbol{x}, \boldsymbol{s}) \quad \text{ for all } \pi \in \mathcal{S}_I \cup \mathcal{S}_{I^c}.$$

Thus Ψ changes its sign if we interchange any particles \mathbf{y}_i and \mathbf{y}_j which possess the same spin, i.e. $s^{(i)} = s^{(j)}$. So the set of particles, and therefore also the set of spacial coordinates, naturally split into two groups I_+ and I_- . In detail, for wavefunctions of d particles \mathbf{y}_i we can (without loss of generality) assume that the first $\#I_+$ indices i belong to the group of positive spin, whereas the rest of the particles have negative spin, that is, $I_+ = \{1, \ldots, \#I_+\}$ and $I_- = I_+^c = \{\#I_+ + 1, \ldots, d\}$.

In physics it is well-known that some problems, e.g., the electronic Schrödinger equation, which involve (general) wavefunctions can be reduced to a bunch of similar problems, where each of them only acts on functions Ψ_s from a certain Hilbert space $\mathcal{F}_d = \mathcal{F}_d(s)$. That is,

$$\Psi_{\boldsymbol{s}} = \Psi(\cdot, \boldsymbol{s}) \in \mathcal{F}_d \subset \{f \colon (\mathbb{R}^3)^d \to \mathbb{R}\}$$

with a given fixed spin configuration $s \in C^d$. Of course every possible spin configuration s corresponds to exactly one choice $I_+ \subseteq \{1,\ldots,d\}$ of indices. Moreover, it is known that \mathcal{F}_d is a Hilbert space with a tensor product structure. Therefore we can model wavefunctions as elements of certain classes of smoothness, e.g., $\mathcal{F}_d \subset H_d = H_1 \otimes \cdots \otimes H_1 = W_2^{(1,\ldots,1)}((\mathbb{R}^3)^d)$, as Yserentant [Yse10] recently did, and incorporate spin properties by using projections of the type $\mathfrak{A} = \mathfrak{A}_{I_+} \circ \mathfrak{A}_{I_-}$, as defined in Section 5.1.1. In particular, Lemma 5.2 then yields

$$\mathcal{F}_d = \mathfrak{A}(H_d) = \mathfrak{A}_{I_+}(H_{\#I_+}) \otimes \mathfrak{A}_{I_-}(H_{\#I_-}),$$

and the system of all

$$\overline{\xi}_{\boldsymbol{k}} = \sqrt{\#\mathcal{S}_{I_{+}} \#\mathcal{S}_{I_{-}}} \, \mathfrak{A}(e_{\boldsymbol{k}}), \quad \boldsymbol{k} \in \overline{\nabla}_{d},$$

with

$$\overline{\nabla}_d = \{ \boldsymbol{k} = (\boldsymbol{i}, \boldsymbol{j}) \in \mathbb{N}^{\#I_+} \times \mathbb{N}^{\#I_-} \mid i_1 < \dots < i_{\#I_+} \text{ and } j_1 < \dots < j_{\#I_-} \}$$

is an orthonormal basis of $\mathcal{F}_d = \mathfrak{A}(H_d)$, where the set $\{e_{\boldsymbol{m}} \mid \boldsymbol{m} \in \mathbb{N}^d\}$ is once again assumed to be an orthonormal tensor product basis of $H_d = H_1 \otimes \cdots \otimes H_1$ constructed with the help of $\{e_m \mid m \in \mathbb{N}\}$, an arbitrary orthonormal basis of H_1 .

Note that in the previous sections the underlying Hilbert space H_1 always consists of univariate functions. In contrast, wavefunctions of one particle depend on at least three (spacial) variables, but we want to stress that this is just a formal issue. Anyway, our approach radically decreases the degrees of freedom and improves the solvability of certain problems $S = (S_d)_{d \in \mathbb{N}}$ like the approximation problem, i.e. $S_d = \mathrm{id} : H_d \to \mathcal{G}_d$ for every $d \in \mathbb{N}$, considered in connection with the electronic Schrödinger equation.

Theorem 5.4 provides an algorithm which is optimal for the \mathcal{G}_d -approximation of d-particle wavefunctions in \mathcal{F}_d with respect to all linear algorithms that use at most n continuous linear functionals. Therefore we only need to choose the right ONB $\{e_m = \phi_m \mid m \in \mathbb{N}\}$ of H_1 which coincides with the eigenfunctions of the univariate operator $W_1 = S_1^{\dagger} S_1$. Moreover, the error can be calculated exactly in terms of the eigenvalues $\lambda = (\lambda_m)_{m \in \mathbb{N}}$ of W_1 .

Furthermore it is possible to prove a modification of Theorem 5.16 for problems dealing with wavefunctions. In fact, for the above mentioned approximation problem, polynomial tractability, as well as strong polynomial tractability, is equivalent to the fact that the sequence λ of the squared singular values of the univariate problem belongs to some ℓ_{τ} -space if we consider the absolute error criterion. The reason is that all the assertions in Section 5.3.3 can be easily extended to the multiple partially antisymmetric case. In detail, if we denote the number of antisymmetric coordinates $\boldsymbol{x}^{(i)}$ within each antisymmetry group $I_d^{(m)} \subseteq \{1,\ldots,d\}$ by $a_{d,m}$ with $m=1,\ldots,M$ then the constraint $a_d+b_d=d$ extends to

$$a_{d,1} + \dots + a_{d,M} + b_d = d.$$

Here b_d again denotes the number of coordinates without any antisymmetry condition. In conclusion, the sufficient condition (5.25) in Proposition 5.12 transfers to

$$\frac{1}{d} \sum_{m=1}^{M} \ln(a_{d,m}!) \ge \|\lambda \mid \ell_{\tau}\|^{\tau} \quad \text{for all } d \ge d_0,$$

which is always satisfied in the case of wavefunctions, since then M=2 and the cardinality $a_{d,m}$ of at least one of the groups of the same spin needs to grow linearly with the dimension d.

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