Distributing many points on spheres: minimal energy and designs

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Abstract

This survey discusses recent developments in the context of spherical designs and minimal energy point configurations on spheres. The recent solution of the long standing problem of the existence of spherical t-designs on $\mathbb{S}^d$ with $\mathcal{O}(t^d)$ number of points by A. Bondarenko, D. Radchenko, and M. Viazovska attracted new interest to this subject. Secondly, D. P. Hardin and E. B. Saff proved that point sets minimising the discrete Riesz energy on $\mathbb{S}^d$ in the hypersingular case are asymptotically uniformly distributed. Both results are of great relevance to the problem of describing the quality of point distributions on $\mathbb{S}^d$, as well as finding point sets, which exhibit good distribution behaviour with respect to various quality measures.

Dedicated to Edward B. Saff on the occasion of his 70th birthday

1. Introduction

The title of this survey is alluding to the fundamental paper by E. B. Saff and A. B. J. Kuijlaars [176], which discusses in large generality how to construct point sets on the sphere

$$\mathbb{S}^d := \left\{ \mathbf{x} \in \mathbb{R}^{d+1} \mid \|\mathbf{x}\| = 1 \right\},$$

which have "good distribution properties" with respect to various measures. That paper appeared in 1997 and motivated and initiated a fruitful direction of research. Evenly distributed point sets have many applications, the most prominent being numerical integration, approximation, interpolation, and sampling; each of these applications needs a different definition of what the word "evenly" should mean.

The present survey on point distributions on the sphere is motivated by two rather recent important contributions, which both shed new light on the subject. The first was D. P. Hardin's and E. B. Saff's [119] proof that minimal energy point configurations on the sphere provide asymptotically uniformly distributed point sets for hypersingular Riesz potentials. This result has become popular under the name *Poppy seed bagel theorem*. This will be one subject of Section 3. The second breakthrough was A. V. Bondarenko's, D. Radchenko's, and M. S. Viazovska's [29] proof that spherical t-designs with $\mathcal{O}(t^d)$ number of points exist on $\mathbb{S}^d$. The definition of spherical designs, their properties, relevance, and occurrence in different contexts will be the subject of Section 2.

Throughout the paper we use the notation $X_N = \{\mathbf{x}_1^{(N)}, \mathbf{x}_2^{(N)}, \ldots, \mathbf{x}_N^{(N)}\}$ for a set of N points; N will always denote the number of points, and the points $\mathbf{x}_1^{(N)}, \mathbf{x}_2^{(N)}, \ldots, \mathbf{x}_N^{(N)}$ usually depend on N. In order to keep the notation simple, we suppress this dependence.

Numerical computation of integrals over multidimensional domains is usually done by weighted sums of point evaluations of the function f to be integrated. Since the sphere is a domain with very high symmetry, it is preferable to use equal weight integration methods,

$$\int_{\mathbb{S}^d} f(\mathbf{x}) \, \mathrm{d}\sigma_d(\mathbf{x}) \approx \frac{1}{N} \sum_{j=1}^{N} f(\mathbf{x}_j). \tag{1.1}$$

Here, and throughout the paper, we denote by σ_d the *normalised* surface area measure on $\mathbb{S}^d$. For non-random collections of integration nodes, using the right-hand side of (1.1) as approximation for the integral is known as a *Quasi-Monte Carlo (QMC) method*. In order to obtain a good approximation of the integral by the sum, the point set X_N should be *well distributed* over the whole sphere $\mathbb{S}^d$. The notion of good distribution of a point set is described in an intrinsic way in this context: the distribution is *good*, if the error for numerical integration is small for a certain set of functions. More precisely, a sequence of node sets (X_N) will be called *asymptotically uniformly distributed*, if the relation

$$\lim_{N \to \infty} \frac{1}{N} \sum_{j=1}^{N} \mathbb{1}_C(\mathbf{x}_j) = \sigma_d(C) \tag{1.2}$$

holds for all spherical caps C (here and throughout the paper $\mathbb{1}_C$ denotes the indicator function of the set C). This is equivalent to

$$\lim_{N \to \infty} \frac{1}{N} \sum_{j=1}^{N} f(\mathbf{x}_j) = \int_{\mathbb{S}^d} f(\mathbf{x}) \, \mathrm{d}\sigma_d(\mathbf{x}) \tag{1.3}$$

for all continuous functions $f : \mathbb{S}^d \to \mathbb{R}$ (see [140]).

Placing N points on the sphere so that the least distance (or equivalently, the smallest angle) between two points,

$$\delta(X_N) := \min_{1 \leq i < j \leq N} \|\mathbf{x}_i - \mathbf{x}_j\| \tag{1.4}$$

is maximised is a further classical approach to obtain evenly distributed points, see [112]. This is the problem of best packing which has attracted attention for a long time in and outside of mathematics. For instance, the best packing problem on the sphere is attributed to P. M. L. Tammes [192], a botanist, who searched for an explanation of the surface structure of pollen grains.

A likewise geometric, but qualitatively different way of evenly distributing N points is to cover the sphere with N spherical caps of radius r centred at these points, and to minimise r. Equivalently, the size of the largest cap, which does not contain a point of X_N,

$$\rho(X_N) := \max_{\mathbf{y} \in \mathbb{S}^d} \min_{1 \le i \le N} \|\mathbf{y} - \mathbf{x}_i\|, \tag{1.5}$$

should be as small as possible. This quantity is called the *covering radius*, and is also known as the *mesh norm* or the *fill radius*. This is the problem of best covering which originates from the realm of facility location problems where the farthest distance of a point on the sphere to the nearest point of X_N (service distance) is minimised.

The definitions of both, the least distance (1.4) and the mesh norm (1.5) can be given using the Euclidean distance or the geodesic distance. Since the geodesic distance on $\mathbb{S}^d$ equals $2 \arcsin(\| \cdot \|/2)$, they are approximately the same for small distances.

Sampling function values on the sphere (*e.g.*, for approximation or interpolation by splines or radial basis functions) or exploring spatial directions in an efficient way requires again *well distributed* (but not necessary uniformly distributed) point sets on $\mathbb{S}^d$. In this context, the quality of a point set is measured differently: well distributed points are required to be *dense* in a quantifiable way. As the mesh norm arises in the error of approximation and good separation is generally associated with the "stability" of an approximation or interpolation method, one would prefer a *quasi-uniform* sequence of point sets (X_N) with uniformly bounded *mesh-separation ratio*

$$\gamma(X_N) := \frac{\rho(X_N)}{\delta(X_N)}. \tag{1.6}$$

The mesh-separation ratio can be regarded as the "condition number" for the point set. Quasi-uniformity is a crucial property for a number of methods (see [99, 144, 180]).

2. Spherical Designs

Spherical designs were initially defined by P. Delsarte, J. M. Goethals, and J. J. Seidel [87] in the context of algebraic combinatorics on spheres. Since then, spherical designs have gained attraction in different areas of mathematics, ranging over number theory, geometry, algebraic and geometric combinatorics, and numerical analysis. We will give an account of these aspects in this section. For a survey on further developments in the context of spherical designs mainly from the point of view of algebraic combinatorics and number theory, we refer to the survey [15] by E. Bannai and E. Bannai.

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2.1. Definition

A spherical t-design is a finite set of points $X_N \subset \mathbb{S}^d$, such that

$$\frac{1}{N} \sum_{\mathbf{x} \in X_N} f(\mathbf{x}) = \int_{\mathbb{S}^d} f(\mathbf{x}) \, d\sigma_d(\mathbf{x}) \tag{2.1}$$

for all polynomials $f \in \mathbb{R}[x_1, \ldots, x_{d+1}]$ (restricted to the sphere $\mathbb{S}^d$) of total degree $\leq t$. This definition is equivalent to

$$\sum_{(\mathbf{x},\mathbf{y}) \in X_N \times X_N} P_\ell^{(d)}(\langle \mathbf{x}, \mathbf{y} \rangle) = 0 \tag{2.2}$$

for $\ell = 1, \ldots, t$; here $P_\ell^{(d)}$ denotes the Legendre polynomial of degree ℓ for the sphere $\mathbb{S}^d$ (see [159]). These polynomials are multiples of Gegenbauer polynomials C_ℓ^α with $\alpha = \frac{d-1}{2}$ normalised so that $P_\ell^{(d)}(1) = 1$. The equivalence of (2.2) and (2.1) is then an immediate consequence of the fact that the restrictions of polynomials to $\mathbb{S}^d$ are spanned by the harmonic polynomials (*i.e.*, polynomials p satisfying $\triangle_{d+1} p = 0$ for the Laplace operator $\triangle_{d+1}$ in $\mathbb{R}^{d+1}$) and the addition theorem for spherical harmonics (see [159])

$$\sum_{k=1}^{Z(d,\ell)} Y_{\ell,k}(\mathbf{x}) Y_{\ell,k}(\mathbf{y}) = Z(d,\ell) \, P_\ell^{(d)}(\langle \mathbf{x}, \mathbf{y} \rangle) \text{ with } Z(d,\ell) := \frac{2\ell + d - 1}{d - 1} \binom{\ell + d - 2}{d - 2}; \tag{2.3}$$

here and throughout this paper $Y_{\ell,k}$ ($k = 1, \ldots, Z(d,\ell)$) denotes a real orthonormal basis of the space of spherical harmonics of total degree ℓ with respect to the scalar product $\langle f, g \rangle := \int_{\mathbb{S}^d} f(\mathbf{x}) g(\mathbf{x}) \, d\sigma_d(\mathbf{x})$. Condition (2.2) is then obtained by considering condition (2.1) for $f = Y_{\ell,k}$, squaring it, summing over k, and using (2.3).

There are two further characterisations of spherical t-designs, which show the connection of this concept to other areas of mathematics. It was observed in [151] that a spherical $2t$-design X_N with N points gives an isometric embedding of ℓ_2^{d+1} into ℓ_{2t}^N, which comes from the identity

$$\frac{1}{N} \sum_{\mathbf{x} \in X_N} \langle \mathbf{x}, \mathbf{a} \rangle^{2t} = \frac{1 \cdot 3 \cdot 5 \cdots (2t - 1)}{(d + 1)(d + 2) \cdots (d + 2t - 1)} \langle \mathbf{a}, \mathbf{a} \rangle^t, \tag{2.4}$$

which is valid for all $\mathbf{a} \in \mathbb{R}^{d+1}$. This identity is an immediate consequence of (2.2) and the expansion of x^{2t} as a sum of Legendre polynomials.

A spherical t-design $X_N = \{\mathbf{x}_1, \mathbf{x}_2, \ldots, \mathbf{x}_N\}$ is called *rigid* (see [14]), if there exists an $\varepsilon > 0$, such that for all t-designs $X_N' = \{\mathbf{x}_1', \mathbf{x}_2', \ldots, \mathbf{x}_N'\}$ with $\|\mathbf{x}_i - \mathbf{x}_i'\| < \varepsilon$ (for $i = 1, \ldots, N$) there exists a rotation $\eta \in \mathrm{SO}(d+1)$ such that $X_N' = \eta X_N$.

It was observed in [107] in passing and later rediscovered in [188] that spherical t-designs can be characterised by a variational property. Let p be a polynomial of degree t given by

$$p(z) = \sum_{\ell=1}^{t} a_\ell P_\ell^{(d)}(z) \quad \text{with } a_\ell > 0 \text{ for } \ell = 1, \ldots, t.$$

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Then X_N is a spherical t-design if and only if the energy functional $E_p(X_N)$ defined by

$$E_p(X_N) := \sum_{(\mathbf{x},\mathbf{y}) \in X_N \times X_N} p(\langle \mathbf{x}, \mathbf{y} \rangle) \tag{2.5}$$

vanishes for X_N. Since the sum (2.5) is non-negative for all X_N, designs are minimisers of this sum. This was used in [188] to characterise designs as stationary points of $E_p(X_N)$.

In order to explain the connection to other questions of algebraic and geometric combinatorics on $\mathbb{S}^d$, we need the following notion. For a finite subset X_N of $\mathbb{S}^d$, we define

$$A(X_N) := \{\langle \mathbf{x}, \mathbf{y} \rangle \mid \mathbf{x}, \mathbf{y} \in X_N, \mathbf{x} \neq \mathbf{y}\}, \tag{2.6}$$

the set of mutual inner products of distinct points. For a given $A \subset [-1, 1)$, X_N is called an A-code, if $A(X_N) \subset A$. Then, for instance, the problem of best packing of N points on $\mathbb{S}^d$ can be formulated as finding the minimal β, such that there exists an A-code with $A = [-1, \beta]$. In particular, the determination of the *kissing number*, that is the maximum number of non-overlapping unit spheres that can touch another given unit sphere, is equivalent to finding the maximal cardinality of an A-code for $A = [-1, \frac{1}{2}]$ (see [160]).

2.2. Relation to lattices

Designs have a very interesting connection to the theory of lattices that we want to explain here. A lattice in $\mathbb{R}^d$ is a $\mathbb{Z}$-module

$$\Lambda := \mathbb{Z}\mathbf{v}_1 \oplus \mathbb{Z}\mathbf{v}_2 \oplus \cdots \oplus \mathbb{Z}\mathbf{v}_d, \tag{2.7}$$

where $\mathbf{v}_1, \ldots, \mathbf{v}_d$ is a basis of $\mathbb{R}^d$. For more detailed information on lattices, we refer to the book [78]. A lattice Λ is called *even*, if all the squared norms $\|\mathbf{v}\|^2$ for $\mathbf{v} \in \Lambda$ are even. Even lattices can only exist if d is divisible by 8.

For any lattice Λ, the dual lattice is defined by

$$\Lambda^{\#} := \{\mathbf{x} \in \mathbb{R}^d \mid \forall \mathbf{v} \in \Lambda : \langle \mathbf{v}, \mathbf{x} \rangle \in \mathbb{Z}\}. \tag{2.8}$$

The set $\Lambda^{\#}$ is again a lattice. A lattice Λ is called *unimodular*, if $\Lambda^{\#} = \Lambda$. For an even unimodular lattice Λ, the ϑ-series

$$\vartheta_\Lambda(\tau) = \sum_{\mathbf{v} \in \Lambda} e^{\pi i \|\mathbf{v}\|^2 \tau}, \qquad \tau \in \mathbb{C}, \quad \mathrm{Im}(\tau) > 0, \tag{2.9}$$

is a modular form of weight $d/2$; *i.e.*, the transformation formula

$$\vartheta_\Lambda(-1/\tau) = \tau^{d/2} \, \vartheta_\Lambda(\tau)$$

holds for all τ with $\mathrm{Im}(\tau) > 0$. This follows from an application of Poisson's summation formula.

The theory of modular forms (see [8]) yields that for a form f of weight $d/2$, given by

$$f(\tau) = \sum_{n=0}^{\infty} a_n \, e^{2\pi i n \tau},$$

at least one of the coefficients a_n for $n = 1, \ldots, 2 + 2\lfloor \frac{d}{24} \rfloor$ has to be non-zero. This implies that every even unimodular lattice Λ contains a non-zero vector $\mathbf{v}$ with $\|\mathbf{v}\| \leq 2 + 2\lfloor \frac{d}{24} \rfloor$. If the shortest non-zero vector $\mathbf{v}$ in Λ satisfies $\|\mathbf{v}\|^2 = 2 + 2\lfloor \frac{d}{24} \rfloor$, then the lattice is called *extremal*.

For a homogeneous polynomial p of degree $j \geq 1$, which is also harmonic ($\triangle_d \, p = 0$), and an even unimodular lattice Λ, the series

$$\vartheta_{\Lambda,p}(\tau) = \sum_{\mathbf{v} \in \Lambda} p(\mathbf{v}) \, e^{\pi i \|\mathbf{v}\|^2 \tau}, \qquad \text{Im}(\tau) > 0, \tag{2.10}$$

is a modular form of weight $d/2 + j$. It is immediate from the symmetry of Λ and the homogeneity of p that $\vartheta_{\Lambda,p} = 0$ if j is odd. If j is even, the series $\vartheta_{\Lambda,p}$ is a cusp form; *i.e.*,

$$\lim_{\text{Im}(\tau) \to \infty} \vartheta_{\Lambda,p}(\tau) = 0,$$

since $p(\mathbf{0}) = 0$.

From the theory of modular forms (see [8]) it is known that there is a unique cusp form $\Delta(\tau)$ (this standard notation "Delta" is not to be confused with the Laplace operator $\triangle$ "triangle") of weight 12 having an expansion of the form

$$\Delta(\tau) = e^{2\pi i \tau} + \sum_{n=2}^{\infty} a_n \, e^{2\pi i n \tau}.$$

It follows from degree considerations that for an even, unimodular, extremal lattice Λ and a homogeneous, harmonic polynomial p of degree $j \geq 1$,

$$\vartheta_{\Lambda,p}(\tau) = \Delta(\tau)^{1 + \lfloor \frac{d}{24} \rfloor} f(\tau), \tag{2.11}$$

where $f(\tau)$ is a modular form of weight

$$\frac{d}{2} - 12 \left\lfloor \frac{d}{24} \right\rfloor + j - 12.$$

If this weight is negative, then f has to vanish identically, since there are no modular forms of negative weight. Consider $d = 24m + k$ with $k \in \{0, 8, 16\}$ and insert this into the above equation. This yields $k/2 + j - 12$ for the weight of f. For $k = 0$, 8, and 16, this yields negative weights for $j \leq 11$, 7, and 3, respectively, which shows that for all shells of the corresponding lattices the sum

$$\sum_{\|\mathbf{v}\|^2 = n} p(\mathbf{v}) = n^{j/2} \sum_{\|\mathbf{v}\|^2 = n} p\left(\mathbf{v}/\sqrt{n}\right) = 0 \tag{2.12}$$

for all homogeneous harmonic polynomials of degree $1 \leq j \leq 11, 7, 3$, respectively. Summing up, we have proved the following theorem.

Theorem 2.1 ([198]). *Let Λ be an extremal even unimodular lattice in $\mathbb{R}^d$ with $d = 24m + k$ for $k = 0, 8, 16$. Then any non-empty shell of Λ defines a j-design $n^{-\frac{1}{2}}\{\mathbf{v} \in \Lambda \mid \|\mathbf{v}\|^2 = n\}$ for $j = 11 - k/2$.*

For further developments in the context of designs, lattices, modular forms, and algebraic codes, we refer to [10, 12, 15, 80, 161, 162].

2.3. Lower bounds

In [87] a linear programming method has been developed, which produces lower bounds for the number of points N of a spherical t-design. This method has been applied successfully to related questions, such as finding the minimal cardinality of $A(X_N)$ for sets X_N with N points.

We give a short explanation of this powerful method, that has been applied with great success to several problems of discrete geometry, the most prominent being the solution of Kepler's conjecture by T. Hales [113]. Since we use the method only on the sphere, we restrict the description to this case here. Let $f : [-1, 1] \to \mathbb{R}^+$ be a continuous positive function given by its expansion in terms of Legendre polynomials; *i.e.,*

$$f(x) = \sum_{n=0}^{\infty} \widehat{f}(n)\, Z(d, n)\, P_n^{(d)}(x). \tag{2.13}$$

If the coefficients $\widehat{f}(n)$ are non-positive for $n > t$, then any t-design X_N has to have cardinality $N \geq f(1)/\widehat{f}(0)$. The proof of this fact is given by the relations

$$f(1)N \leq \sum_{(\mathbf{x},\mathbf{y}) \in X_N \times X_N} f(\langle \mathbf{x}, \mathbf{y} \rangle)$$

$$= \widehat{f}(0)\, (N)^2 + \sum_{n > t} \widehat{f}(n)\, Z(d, n) \sum_{(\mathbf{x},\mathbf{y}) \in X_N \times X_N} P_n^{(d)}(\langle \mathbf{x}, \mathbf{y} \rangle)$$

$$\leq \widehat{f}(0)N^2.$$

The first inequality is a consequence of the positivity of f, the equality uses the characterisation of t-designs by (2.2), and the second inequality uses that $\widehat{f}(n) \leq 0$ for $n > t$ and the non-negativity of the double sums $\sum_{(\mathbf{x},\mathbf{y}) \in X_N \times X_N} P_n^{(d)}(\langle \mathbf{x}, \mathbf{y} \rangle)$. Equality can only occur, if $f(x) = 0$ for all $x \in A(X_N)$ and $\widehat{f}(n) = 0$ for $n > t$; *i.e.,* f is a polynomial of degree $\leq t$.

In [87] a polynomial p_t of degree t was constructed, which gives the lower bound

$$N \geq \begin{cases} \binom{d+t/2}{d} + \binom{d+t/2-1}{d} & \text{for } t \text{ even,} \\ 2\binom{d+(t-1)/2}{d} & \text{for } t \text{ odd.} \end{cases} \tag{2.14}$$

This polynomial actually provides the best possible lower bound that can be obtained by polynomial functions f. Designs attaining this lower bound are called *tight*. It was shown in [16, 17] that tight designs only exist for $d = 1$ and all t, or finitely many values of t if

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$d \geq 2$. The proof uses the fact that for a tight spherical t-design X_N, the corresponding set $A(X_N)$ has to consist exactly of the zeros of the polynomial p_t constructed in [87]. Galois theory is then used to derive a contradiction, if t is large enough. The same fact about $A(X_N)$ shows that tight designs are rigid (see [14]). Notice that the bound is of order t^d. Later, V. A. Yudin [202] considered a wider class of functions than polynomials. He constructed a function f, which allowed for a considerable improvement of the lower bound. The bounds for spherical designs he gives are again of order t^d, but they are larger by a factor depending on d. This factor is exponential in d. For $d = 2$, the gain is asymptotically about 7%.

We give a short description of Yudin's construction. The function f is obtained as the spherical convolution of two positive functions F and G, which ensures the positivity of f. The spherical convolution of two functions F and G is given by

$$F \star G(\langle \mathbf{x}, \mathbf{y} \rangle) := \int_{\mathbb{S}^d} F(\langle \mathbf{x}, \mathbf{z} \rangle) G(\langle \mathbf{z}, \mathbf{y} \rangle) \, d\sigma_d(\mathbf{z}).$$

In order to obtain the required sign change of the Laplace-Fourier coefficients of f, the function G is chosen as

$$G(x) = (1 - x^2)^{1 - \frac{d}{2}} \left((1 - x^2)^{\frac{d}{2}} F'(x) \right)' + (t + 1)(t + d) F(x),$$

which is motivated by the fact that the Legendre polynomials are eigenfunctions of the differential operator L defined by

$$LF(x) := (1 - x^2)^{1 - \frac{d}{2}} \left((1 - x^2)^{\frac{d}{2}} F'(x) \right)'.$$

If F is expressed in terms of its Laplace-Fourier expansion

$$F(x) = \sum_{n=0}^{\infty} \widehat{F}(n) \, Z(d, n) \, P_n^{(d)}(x),$$

then the function G is given by

$$G(x) = \sum_{n=0}^{\infty} (t + 1 - n)(n + t + d) \, \widehat{F}(n) \, Z(d, n) \, P_n^{(d)}(x).$$

Hence the spherical convolution is given as

$$f(x) = F \star G(x) = \sum_{n=0}^{\infty} (t + 1 - n)(n + t + d) \, \widehat{F}(n)^2 \, Z(d, n) \, P_n^{(d)}(x),$$

which has the required sign-change in its Laplace-Fourier coefficients. The problem of maximising the quotient $f(1)/\widehat{f}(0) = f(1)/\widehat{F}(0)^2$ turns out to be a variational problem for

F:

$$f(1) = \frac{\omega_{d-1}}{\omega_d} \int_{-1}^{1} F(x)G(x) \left(1 - x^2\right)^{d/2-1} \, \mathrm{d}x$$

$$= \frac{\omega_{d-1}}{\omega_d} \int_{-1}^{1} \left(- \left(1 - x^2\right) F'(x)^2 + (t+1)(t+d) F(x)^2\right) \left(1 - x^2\right)^{d/2-1} \, \mathrm{d}x$$

must be maximised subject to the condition

$$\widehat{F}(0) = \frac{\omega_{d-1}}{\omega_d} \int_{-1}^{1} F(x) \left(1 - x^2\right)^{d/2-1} \, \mathrm{d}x = 1.$$

The symbol ω_d denotes the surface area of $\mathbb{S}^d$ and it satisfies

$$\int_{-1}^{1} \left(1 - x^2\right)^{d/2-1} \, \mathrm{d}x = \frac{\omega_d}{\omega_{d-1}} = \frac{\sqrt{\pi}\,\Gamma(d/2)}{\Gamma((d+1)/2)}.$$

Furthermore, f has to be non-negative on $[-1, 1]$. This is achieved by assuming that F and G are both non-negative. The solution of this variational problem with the additional condition on the sign of F is then given by

$$F(x) = \begin{cases} P_{t+1}^{(d)}(x) - P_{t+1}^{(d)}(\alpha_t) & \text{for } \alpha_t \leq x \leq 1, \\ 0 & \text{for } -1 \leq x \leq \alpha_t, \end{cases}$$

where α_t is the largest zero of $\frac{d}{dx} P_{t+1}^{(d)}(x)$. In this case G is a piecewise constant function. Putting everything together yields the lower bound

$$N \geq \frac{\int_{-1}^{1} \left(1 - x^2\right)^{d/2-1} \, \mathrm{d}x}{\int_{\alpha_t}^{1} \left(1 - x^2\right)^{d/2-1} \, \mathrm{d}x} = \frac{1}{\sigma_d(\{\mathbf{y} \in \mathbb{S}^d \mid \langle \mathbf{y}, \mathbf{x} \rangle \geq \alpha_t\})} \gg_d t^d \tag{2.15}$$

for a t-design X_N. For a more detailed exposition, we refer to [202]. Even if the obtained function f gives a better value for the lower bound than the polynomial given in [87], the technical requirement of non-negativity of F seems to leave room for further improvement.

A similar construction for a function f is used in [73] to obtain linear programming bounds for the packing density of spheres in $\mathbb{R}^d$. The function obtained there has similar features; in particular, it is supported on a short interval. In [73, Section 5] it is mentioned that the function constructed by this convolution method does **not** produce optimal bounds (see [72]). The reason seems to be exactly the non-negativity requirement on the corresponding function F. Thus there is reason to believe that Yudin's lower bound for the cardinality of t-designs can still be improved.

2.4. Existence

On the other hand, the question of the existence of spherical t-designs has been answered affirmatively. First, a rather general result obtained in [184] shows that, for given t and

large enough N, there exists a t-design with N points. Actually, the result given in [184] is more general: given a path connected topological space Ω and a finite measure μ that charges every non-empty open set, then for any finite set of continuous real valued functions $f_1, \ldots, f_t$ there exists an N_0, such that for every $N > N_0$ there exists a set $X_N \subset \Omega$, such that

$$\frac{1}{N} \sum_{x \in X_N} f_j(x) = \frac{1}{\mu(\Omega)} \int_\Omega f_j(x) \, \mathrm{d}\mu(x)$$

for $j = 1, \ldots, t$. The result gives a bound for the number of points needed by geometric quantities defined in terms of Ω and the functions $f_1, \ldots, f_t$. These quantities are very difficult to compute even in the special case of the sphere.

In order to discuss the question of existence of spherical designs further, we introduce two quantities:

$$N(t, d) := \min \left\{ N \mid \exists X_N \subset \mathbb{S}^d : X_N \text{ is a } t\text{-design} \right\},$$
$$N^*(t, d) := \min \left\{ K \mid \forall N \geq K : \exists X_N \subset \mathbb{S}^d : X_N \text{ is a } t\text{-design} \right\}.$$

By definition it is clear that $N^*(t, d) \geq N(t, d)$.

Only very recently, the long standing problem of the existence of spherical t-designs with $\mathcal{O}(t^d)$ points was answered affirmatively by A. V. Bondarenko, D. Radchenko, and M. S. Viazovska [29]. They proved that $N^*(t, d) = \mathcal{O}(t^d)$ with an explicit, but large implied constant. The proof puts the existence of spherical t-designs with $N = \mathcal{O}(t^d)$ in the context of Brouwer's degree theorem: Let $\mathcal{P}_t$ denote the Hilbert space of polynomials P of total degree $\leq t$ with

$$\int_{\mathbb{S}^d} P(\mathbf{x}) \, \mathrm{d}\sigma_d(\mathbf{x}) = 0$$

equipped with the scalar product

$$\langle P, Q \rangle = \int_{\mathbb{S}^d} P(\mathbf{x}) Q(\mathbf{x}) \, \mathrm{d}\sigma_d(\mathbf{x}).$$

Then for every $\mathbf{x} \in \mathbb{S}^d$ there exists a polynomial $G_{\mathbf{x}}$ such that

$$\langle G_{\mathbf{x}}, Q \rangle = Q(\mathbf{x})$$

for all $Q \in \mathcal{P}_t$. A set of points X_N then is a spherical t-design, if

$$G_{\mathbf{x}_1} + \cdots + G_{\mathbf{x}_N} \equiv 0.$$

Then continuous maps $\mathbf{x}_i : \mathcal{P}_t \to \mathbb{S}^d$ are constructed, which in turn define the map $f : \mathcal{P}_t \to \mathcal{P}_t$ by

$$f(P) := G_{\mathbf{x}_1(P)} + \cdots + G_{\mathbf{x}_N(P)}.$$

Then for any $P \in \mathcal{P}_t$ the identity

$$\langle P, f(P) \rangle = \sum_{i=1}^{N} P(\mathbf{x}_i(P))$$

holds. A polynomial $\tilde{P}$ with $f(\tilde{P}) = 0$ then gives a spherical t-design $\{\mathbf{x}_1(\tilde{P}), \ldots, \mathbf{x}_N(\tilde{P})\}$. The construction of the maps $\mathbf{x}_i$ is the crucial part of the proof. Starting with points $\{\mathbf{x}_1(0), \ldots, \mathbf{x}_N(0)\}$ in the parts of an equal area partition of the sphere, the maps $\mathbf{x}_i$ are defined by an intricate geometric procedure. The proof finishes by considering the set

$$\Omega = \left\{ P \in \mathcal{P}_t \mid \int_{\mathbb{S}^d} \|\nabla P(\mathbf{x})\| \, \mathrm{d}\sigma_d(\mathbf{x}) < 1 \right\}$$

and observing that the construction of $\mathbf{x}_i$ yields

$$\langle P, f(P) \rangle > 0$$

for $P \in \partial\Omega$. This last inequality is verified by an application of a spherical version of the Marcinkiewicz–Zygmund inequality. In this step the precise choice of the number N of points in relation to t is significant. The Brouwer degree theorem gives the existence of a point $\tilde{P}$ with $f(\tilde{P}) = 0$, which the yields the desired spherical t-design $\{\mathbf{x}_1(\tilde{P}), \ldots, \mathbf{x}_N(\tilde{P})\}$.

In a recent paper [30], the same authors showed that there exist well separated t-designs with optimal order of the number of points:

Theorem 2.2. *Let $d \geq 2$. Then there exist positive constants C_d and β_d such that for every $N \geq C_d t^d$ there exists a t-design X_N and*

$$\|\mathbf{x} - \mathbf{y}\| \geq \beta_d \, N^{-1/d} \qquad \text{for all } \mathbf{x}, \mathbf{y} \in X_N \text{ with } \mathbf{x} \neq \mathbf{y}.$$

The proof is a refinement of their original proof by keeping control on the distance of distinct points.

2.5. Numerical results

Besides the theoretical investigation of the existence of spherical designs, several attempts were made to compute lists of t-designs for moderately large values of t and N. Since most of these computations have been done for $d = 2$, we will restrict to that case in this section.

A first list of t-designs for $t \leq 21$ was provided by R. H. Hardin and N. J. A. Sloane (see [123, 124, 125]). Their list is still available on the web [126]. These numerical computations, as well as those performed in [188], seem to suggest that $N(t, 2)$ is close to $\frac{1}{2}(t+1)^2$. Furthermore, X. Chen and R. S. Womersley found spherical t-designs by numerical computations for $t \leq 100$ (see [68]). Their computations seem to indicate that there exist t-designs with less than $(t+1)^2$ points. Later X. Chen, A. Frommer, and B. Lang [67] used interval arithmetic to prove that there exist spherical t-designs with $(t+1)^2$ points for $1 \leq t \leq 100$. Recently, M. Gräf and D. Potts [108, 109] derived a new method based on fast Fourier transform, which allows to find t-designs numerically with high precision for values of t up to 1000. They also provide their results on the web [110]. Again these numerical results show that $N(t, 2)$ is close to $\frac{1}{2}(t+1)^2$ for small values of t.

Table 1: The lower bounds for the number of points of a t-design derived in [87] (DGS(t)) and in [202] (Y(t)) compared to the number of points of t-designs $N(t,2)$ obtained in [109] (as provided on the website [110]).

t	5	7	9	10	20	30	40	50	60	70	80	90	100	114	124
DGS(t)	12	20	30	36	121	256	441	676	961	1296	1681	2116	2601	3364	3969
Y(t)	12	20	31	37	127	271	470	723	1031	1394	1810	2282	2808	3635	4292
$N(t,2)$	12	24	48	60	216	480	840	1296	1860	2520	3276	4140	5100	6612	7812
$\lfloor \frac{(t+1)^2}{2} \rfloor$	18	32	50	60	220	480	840	1300	1860	2520	3280	4140	5100	6612	7812

2.6. Designs and uniform distribution

It was observed independently in [107] and in [136] that spherical designs provide well-distributed point sets on the sphere.

The *spherical cap discrepancy* of a point set X_N of N points is given by

$$D(X_N) := \sup_{\substack{\mathbf{y} \in \mathbb{S}^d, \\ \varphi \in [0,\pi]}} \left| \frac{1}{N} \sum_{j=1}^{N} \mathbb{1}_{C(\mathbf{y},\varphi)}(\mathbf{x}_j) - \sigma_d(C(\mathbf{y},\varphi)) \right|; \qquad (2.16)$$

the supremum is extended over all spherical caps

$$C(\mathbf{y},\varphi) := \left\{ \mathbf{x} \in \mathbb{S}^d \mid \langle \mathbf{x}, \mathbf{y} \rangle > \cos(\varphi) \right\},$$

and measures the maximum deviation between the empirical distribution of the point set X_N from uniform distribution. In [105] the estimate

$$D(X_N) \leq \frac{C_1(d)}{M} + \sum_{\ell=1}^{M} \frac{C_2(d)}{\ell} \sum_{k=1}^{Z(d,\ell)} \left| \frac{1}{N} \sum_{j=1}^{N} Y_{\ell,k}(\mathbf{x}_j) \right| \qquad (2.17)$$

was proved; here M is an arbitrary positive integer, and $C_1(d)$ and $C_2(d)$ are explicit constants depending only on the dimension d. A similar inequality was later given in [147]. The inequality (2.17) resembles the classical Erdős-Turán-Koksma inequality estimating the Euclidean discrepancy of a point set in $[0,1]^d$ in terms of trigonometric sums (see [140]). In [138], N. M. Korobov introduced *good lattice points* $(g_1, \ldots, g_d) \in \mathbb{Z}^d$ by the requirement that the point set $\{(\{\frac{jg_1}{N}\}, \ldots, \{\frac{jg_d}{N}\}) \mid j = 0, \ldots, N-1\}$ has small discrepancy. Here, $\{x\}$ denotes the fractional part of x, which puts the point set into the unit cube $[0,1]^d$.

In [107] spherical t-designs were regarded as spherical analogues of good lattice points, in the sense that the estimate (2.17) becomes particularly simple when applied to a t-design and choosing the parameter M to be t: the estimate then reduces to

$$D(X_N) \leq \frac{C_1(d)}{t}.$$

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Similarly, for a continuous function $f : \mathbb{S}^d \to \mathbb{R}$ satisfying the Lipschitz-condition $|f(\mathbf{x}) - f(\mathbf{y})| \le C_f \arccos(\langle \mathbf{x}, \mathbf{y} \rangle)$, the estimate

$$\left| \frac{1}{N} \sum_{\mathbf{x} \in X_N} f(\mathbf{x}) - \int_{\mathbb{S}^d} f(\mathbf{x}) \, d\sigma_d(\mathbf{x}) \right| \le C_f \left(6 \frac{d}{M} + \pi \sum_{\ell=1}^{2M} \sum_{k=1}^{Z(d,\ell)} \left| \frac{1}{N} \sum_{\mathbf{x} \in X_N} Y_{\ell,k}(\mathbf{x}) \right| \right) \quad (2.18)$$

was shown in the same paper. Again, taking X_N to be a $2t$-design and $M = t$ gives an estimate $6dC_f/t$ for the integration error.

J. Korevaar and J. L. H. Meyers [136] take a potential theoretic point of view. In their papers [136, 137] they conjecture that $N(t, d) = \mathcal{O}(t^d)$, which was finally proved in [29]. Let $\mu_N := \frac{1}{N} \sum_{\mathbf{x} \in X_N} \delta_{\mathbf{x}}$ be the discrete equal weight distribution supported on $X_N \subset \mathbb{S}^d$. Then the deviation of μ_N and the equilibrium measure σ_d is measured by the deviation of the potential

$$U_{d-1}^{\mu_N}(\mathbf{x}) = \frac{1}{N} \sum_{\mathbf{y} \in X_N} \frac{1}{\|\mathbf{x} - \mathbf{y}\|^{d-1}} = \int_{\mathbb{S}^d} \frac{1}{\|\mathbf{x} - \mathbf{y}\|^{d-1}} \, d\mu_N(\mathbf{y}) \quad (2.19)$$

from the equilibrium potential

$$U_{d-1}^{\sigma_d}(\mathbf{x}) = \int_{\mathbb{S}^d} \frac{1}{\|\mathbf{x} - \mathbf{y}\|^{d-1}} \, d\sigma_d(\mathbf{y}) = 1$$

for $\|\mathbf{x}\| \le r < 1$.

Taking a spherical t-design X_N and using the identity (see [152])

$$\frac{1}{\|r\mathbf{x} - \mathbf{y}\|^{d-1}} = \sum_{n=0}^{\infty} \binom{n+d-2}{d-2} P_n^{(d)}(\langle \mathbf{x}, \mathbf{y} \rangle) \, r^n,$$

we obtain

$$U_{d-1}^{\mu_N}(r\mathbf{x}) - 1 = \sum_{n=t+1}^{\infty} \binom{n+d-2}{d-2} r^n \frac{1}{N} \sum_{\mathbf{y} \in X_N} P_n^{(d)}(\langle \mathbf{x}, \mathbf{y} \rangle). \quad (2.20)$$

Estimating the right hand side by $|P_n^{(d)}(\cdot)| \le 1$, we obtain

$$\left| U_{d-1}^{\mu_N}(r\mathbf{x}) - 1 \right| \le \sum_{n=t+1}^{\infty} \binom{n+d-2}{d-2} r^n. \quad (2.21)$$

The sum on the right hand side can be expressed in closed form.

$$\sum_{n=t+1}^{\infty} \binom{n+d-2}{d-2} r^n$$

$$= (d-1) \binom{t+d-1}{d-1} \frac{1}{(1-r)^{d-1}} \int_0^r (1-\rho)^{d-2} \rho^t \, d\rho$$

$$= \frac{r^{t+1}}{(1-r)^{d-1}} (d-1) \binom{t+d-1}{d-1} \sum_{\ell=0}^{d-2} \binom{d-2}{\ell} (-1)^\ell \frac{r^\ell}{t+\ell+1},$$

which can be proved by multiplying with $(1-r)^{d-1}$ and differentiating. Estimating the integral in the second line by $r^{t+1}/(t+1)$ we get the estimate

$$\left|U_{d-1}^{\mu_N}(r\mathbf{x}) - 1\right| \leq \binom{t+d-1}{d-2} \frac{r^{t+1}}{(1-r)^{d-1}}, \tag{2.22}$$

valid for any t-design X_N, $\mathbf{x} \in \mathbb{S}^d$, and $0 \leq r < 1$. This, together with the fact that $N(t,d) = \mathcal{O}(t^d)$, generalises [136, Theorem 2.2] to arbitrary dimension.

2.7. Applications to numerical integration on $\mathbb{S}^d$

Equal weight quadrature formulas based on spherical designs are exact on polynomials up to a given degree. This puts them into the general framework of QMC integration methods. It turns out that designs are especially useful for integrating functions taken from suitably defined Sobolev spaces on $\mathbb{S}^d$. These spaces are *reproducing kernel Hilbert spaces*, which makes the study of the *worst case error* in integration particularly simple and transparent (see [165]).

In order to describe the results on numerical integration in more detail, we give a precise definition of the function spaces. The negative Laplace-Beltrami operator $-\triangle_d^*$ on $\mathbb{S}^d$ has the eigenvalues $\lambda_\ell := \ell(\ell + d - 1)$, $\ell \in \mathbb{N}_0$. The space of eigenfunctions for the eigenvalue λ_ℓ is spanned by the real spherical harmonics $Y_{\ell,k}$ for $k = 1, \ldots, Z(d,\ell)$. Since the eigenfunctions of $-\triangle_d^*$ form a complete L^2-orthogonal system, every function $f \in L^2(\mathbb{S}^d)$ can be represented by its Laplace-Fourier series expansion

$$f(\mathbf{x}) = \sum_{\ell=0}^{\infty} \sum_{k=1}^{Z(d,\ell)} \widehat{f}_{\ell,k} \, Y_{\ell,k}(\mathbf{x}), \tag{2.23}$$

where the Laplace-Fourier coefficients are given by

$$\widehat{f}_{\ell,k} = \int_{\mathbb{S}^d} f(\mathbf{x}) \, Y_{\ell,k}(\mathbf{x}) \, \mathrm{d}\sigma_d(\mathbf{x}).$$

The series (2.23) has to be interpreted in the L^2-sense. Furthermore, Parseval's identity

$$\sum_{\ell=0}^{\infty} \sum_{k=1}^{Z(d,\ell)} \left|\widehat{f}_{\ell,k}\right|^2 = \int_{\mathbb{S}^d} |f(\mathbf{x})|^2 \, \mathrm{d}\sigma_d(\mathbf{x}) = \|f\|_2^2 \tag{2.24}$$

and the Funk-Hecke formula

$$\sum_{k=1}^{Z(d,\ell)} \widehat{f}_{\ell,k} \, Y_{\ell,k}(\mathbf{x}) = Z(d,\ell) \int_{\mathbb{S}^d} f(\mathbf{y}) \, P_\ell^{(d)}\big(\langle \mathbf{x}, \mathbf{y} \rangle\big) \, \mathrm{d}\sigma_d(\mathbf{y}) \tag{2.25}$$

hold. For more details on the harmonic analysis on $\mathbb{S}^d$, we refer to [159].

In the following we adopt the notation of [61]. For $s \geq 0$, we define the Hilbert space

$$H^s(\mathbb{S}^d) := \left\{ f \in L^2(\mathbb{S}^d) \,\,\Big|\,\, \sum_{\ell=0}^{\infty}(1+\ell)^{2s} \sum_{k=1}^{Z(d,\ell)} \left|\widehat{f}_{\ell,k}\right|^2 < \infty \right\}. \tag{2.26}$$

The inner product on $H^s(\mathbb{S}^d)$ is then given by

$$\langle f, g \rangle_{H^s} := \sum_{\ell=0}^{\infty}(1+\ell)^{2s} \sum_{k=1}^{Z(d,\ell)} \widehat{f}_{\ell,k}\overline{\widehat{g}_{\ell,k}} \tag{2.27}$$

with the corresponding norm $\|f\|_{H^s} := \langle f, f \rangle_{H^s}^{1/2}$. We remark that the "sequence of weights" $((1+\ell)^{2s})$ can be replaced by any other *comparable* sequence (w_ℓ), in the sense of

$$\exists C_1, C_2 > 0 : \forall \ell \in \mathbb{N}_0 : C_1(1+\ell)^{2s} \leq w_\ell \leq C_2(1+\ell)^{2s}.$$

From the definition it is clear that the spaces $H^s(\mathbb{S}^d)$ are getting smaller as the index of smoothness s increases. Furthermore, the Sobolev embedding theorem ensures that $H^s(\mathbb{S}^d)$ embeds continuously into $C^k(\mathbb{S}^d)$ when $s > k + d/2$; in particular, $H^s(\mathbb{S}^d)$ embeds continuously into $C(\mathbb{S}^d)$ when $s > d/2$.

As a consequence of the continuous embedding of $H^s(\mathbb{S}^d)$ into $C(\mathbb{S}^d)$ for $s > d/2$, point evaluation of a function is a continuous functional, which can be represented as a scalar product by the Riesz representation theorem. This ensures the existence of a reproducing kernel given by

$$\mathrm{K}^{(s)}\left(\langle \mathbf{x}, \mathbf{y} \rangle\right) := \sum_{\ell=0}^{\infty} (1+\ell)^{-2s}\, Z(d,\ell)\, P_\ell^{(d)}\left(\langle \mathbf{x}, \mathbf{y} \rangle\right). \tag{2.28}$$

It can be verified immidiately that the function $\mathrm{K}^{(s)}$ is indeed a reproducing kernel satisfying

$$\forall \mathbf{y} \in \mathbb{S}^d : \mathrm{K}^{(s)}(\langle \cdot, \mathbf{y} \rangle) \in H^s(\mathbb{S}^d) \tag{2.29}$$

and

$$f(\mathbf{x}) = \langle f, \mathrm{K}^{(s)}(\langle \cdot, \mathbf{x} \rangle) \rangle_{H^s} \tag{2.30}$$

for all $f \in H^s(\mathbb{S}^d)$ and for all $\mathbf{x} \in \mathbb{S}^d$.

For $f \in H^s(\mathbb{S}^d)$, $s > d/2$, the integration error of the QMC method with node set X_N is given by

$$\mathrm{Error}(X_N)(f) := \frac{1}{N} \sum_{\mathbf{x} \in X_N} f(\mathbf{x}) - \int_{\mathbb{S}^d} f(\mathbf{x})\, \mathrm{d}\sigma_d(\mathbf{x}) = \langle f, R(X_N) \rangle_{H^s}, \tag{2.31}$$

where the function

$$R(X_N)(\mathbf{y}) := \frac{1}{N} \sum_{\mathbf{x} \in X_N} \mathrm{K}^{(s)}(\langle \mathbf{y}, \mathbf{x} \rangle) - 1 \tag{2.32}$$

is called the *representer of the integration error*. It is now a consequence of elementary Hilbert space theory that the worst case error takes the form

$$\mathrm{wce}_{H^s}(X_N) := \sup_{\|f\|_{H^s}=1} |\mathrm{Error}(X_N)(f)| = \|R(X_N)\|_{H^s}. \qquad (2.33)$$

The squared norm $\|R(X_N)\|^2 = \langle R(X_N), R(X_N) \rangle$ can be expressed in terms of the kernel function $\mathrm{K}^{(s)}$ by means of

$$\|R(X_N)\|^2_{H^s} = \frac{1}{N^2} \sum_{\mathbf{x},\mathbf{y}\in X_N} \mathrm{K}^{(s)}(\langle \mathbf{x}, \mathbf{y} \rangle) - 1. \qquad (2.34)$$

For a more precise explanation of this formalism we refer to [61]. The expression (2.34) is a special case of an energy functional as discussed further in Section 3.

Taking a spherical t-design for the set X_N in (2.34), the first t terms in the Laplace-Fourier expansion of $\mathrm{K}^{(s)}$ are annihilated:

$$\mathrm{wce}_{H^s}(X_N)^2 = \sum_{\ell=t+1}^{\infty} (1+\ell)^{-2s} \, Z(d,\ell) \frac{1}{N^2} \sum_{\mathbf{x},\mathbf{y}\in X_N} P_\ell^{(d)}(\langle \mathbf{x}, \mathbf{y} \rangle). \qquad (2.35)$$

Trivially estimating this sum and taking square roots would give a bound $\mathcal{O}(t^{d/2-s})$ for the worst case error.

A method is introduced in [131] and [132], which provides much better estimates for the worst case error in $H^s(\mathbb{S}^2)$. This method was extended and generalised to higher dimensions in [60]. This is now a standard technique in this context. While too technical to give a precise description here, we explain the two main steps. For a full description of the method we refer to [60].

First, the tail of the series defining the kernel

$$\mathrm{K}^{(s)}_t(x) := \sum_{\ell=t+1}^{\infty} (1+\ell)^{-2s} \, Z(d,\ell) \, P_\ell^{(d)}(x)$$

is rewritten in terms of Jacobi polynomials instead of Legendre polynomials. This is done using the fact that the Legendre polynomials are special cases of Jacobi polynomials and that there exist connection formulas involving hypergeometric expressions (see [152]). After this transformation, the function $K^{(s)}_t$ is expressed as a polynomial of degree t plus a series $\widetilde{K}^{(s)}_t$ involving higher order polynomials ("kernel splitting method"). The polynomial part is integrated exactly by the QMC method supported in X_N; the remaining part

$$\sum_{\mathbf{x},\mathbf{y}\in X_N} \widetilde{\mathrm{K}}^{(s)}_t(\langle \mathbf{x}, \mathbf{y} \rangle)$$

is estimated using bounds for Jacobi polynomials. This part is rather delicate, since different estimates have to be used for $\langle \mathbf{x}, \mathbf{y} \rangle \in [-1+c/t, 1-c/t]$ and $|\langle \mathbf{x}, \mathbf{y} \rangle| > 1 - c/t$. Here,

a coarse equidistribution property of spherical t-designs is used to obtain optimal order estimates: there exist constants c_1 and c_2, independent of t and N, such that

$$\# \left(X_N \cap C \left(\mathbf{x}, \frac{c_1}{t} \right) \right) \leq c_2 N \sigma_d \left(C \left(\mathbf{x}, \frac{c_1}{t} \right) \right). \tag{2.36}$$

It was proved in [171] (see also [172]) that this is a general property of positive weight quadrature formulas, which integrate polynomials of degree $\leq t$ exactly; thus (2.36) holds *a forteriori* for spherical t-designs.

Using the technique described above, it was proved in [60] that the worst case integration error on $H^s(\mathbb{S}^d)$ (for $s > d/2$) is

$$\mathrm{wce}_{H^s}(X_N) = \mathcal{O}_{s,d}(t^{-s}). \tag{2.37}$$

Using the recent progress on the existence of spherical designs with optimal growth order (see [29]), this gives

$$\mathrm{wce}_{H^s}(X_N) = \mathcal{O}_{s,d}(N^{-s/d}) \tag{2.38}$$

if X_N is a spherical t-design with $N = \mathcal{O}(t^d)$. This is known to be the optimal order for the worst case error in $H^s(\mathbb{S}^d)$ (see [129, 130]).

The estimate (2.38) led to the following definition in [61]. Let $s > d/2$, then a sequence of point sets (X_N) on $\mathbb{S}^d$ is called a *sequence of QMC-designs* for $H^s(\mathbb{S}^d)$ if (2.38) holds. The supremum over all s for which (2.38) holds, is called the *strength* of the sequence (X_N). By the above description, a sequence of spherical t-designs of optimal growth order $\mathcal{O}(t^d)$ has strength ∞. For a more detailed exposition of QMC-designs and their properties, we refer to the original paper [61]. These investigations were extended to the L^p-setting in [51].

2.8. Applications in approximation and interpolation

Well distributed point sets can also be used as sample points for interpolation formulas. Let $f : \mathbb{S}^d \to \mathbb{R}$ be a continuous function and X_N a set of points with $N \geq Z(d+1, t)$. Observe that

$$Z(d+1, t) = \sum_{k=0}^{t} Z(d, k)$$

is the dimension of the space of all polynomials of total degree $\leq t$ on $\mathbb{S}^d$. Then we seek a polynomial p of degree $\leq t$, such that

$$p(\mathbf{x}_i) = f(\mathbf{x}_i) \qquad \text{holds for } i = 1, \dots, N. \tag{2.39}$$

Choosing $Y_{\ell,k}$ ($\ell = 0, \dots, t$, $k = 1, \dots, Z(d, \ell)$) as a basis for the polynomials and $N = Z(d+1, t)$, we have to solve a system of N linear equations; the points must be chosen so that this system has full rank. One possible method for choosing the set X_N is to maximise the modulus of the determinant of the system, which optimises the numerical stability of solving the system. This approach is in the spirit of finding the extremal

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value of a functional depending on the points, such as the energy functionals discussed in Section 3. This has been done in [187] for $d = 2$.

In [6] it is proposed to use spherical t-designs as interpolation points; for $N \geq Z(d+1, t)$ the determinant of the matrix

$$H_t := \left(\sum_{j=1}^{N} Y_{\ell,k}(\mathbf{x}_j) Y_{\ell',k'}(\mathbf{x}_j) \right)_{(\ell,k),(\ell',k')} \qquad 1 \leq \ell, \ell' \leq t, \quad 1 \leq k \leq Z(d,\ell), \quad 1 \leq k' \leq Z(d,\ell'),$$

is maximised under the constraint that the point set X_N is a spherical t-design. This leads to the definition of *well conditioned* designs. In [6] this is worked out for $d = 2$.

It should be mentioned that, based on the numerical experiments in [110, 126], there exist t-designs with $\lfloor (t+1)^2/2 \rfloor$ points for $t \leq 100$. In the context of interpolation, the number of points has to be chosen $\geq (t+1)^2$, which gives further freedom for choosing the points that can be used to maximise the determinant of H_t.

In [7], the application of spherical designs in the context of approximation of functions $f : \mathbb{S}^d \to \mathbb{R}$ is proposed. Usually, the approximation is computed by integrating f against an approximation kernel G_L, which is a polynomial of degree L in $\langle \mathbf{x}, \mathbf{y} \rangle$ (see [23, 173]). In [7] the integral is replaced by the equal weight quadrature rule given by a $2L$-design X_N. This approach is worked out for $d = 2$ and various classical approximation kernels. In order to make this procedure numerically more stable, a regularisation procedure is investigated: the polynomial p is chosen to minimise

$$\sum_{j=1}^{N} (p(\mathbf{x}_j) - f(\mathbf{x}_j))^2 + \lambda \sum_{j=1}^{N} (\mathcal{R}p(\mathbf{x}_j))^2 \qquad (2.40)$$

amongst all polynomials of degree L; here $\mathcal{R}$ is an operator on the space of polynomials and $\lambda > 0$ is the regularisation parameter. The fact that X_N is chosen as a $2L$-design simplifies the linear algebra behind the least square approximation.

3. Energy

3.1. Minimal energy in applications

A surprising number of diverse applications can be formulated as a discrete or continuous minimal energy problem or a mixture of both. In the discrete setting this means finding a collection of N distinct points in a subset $\Omega \subset \mathbb{R}^p$, $p \geq 1$, that minimises a *discrete* K-*energy functional*

$$E_{\mathrm{K,Q}}(\mathbf{x}_1, \ldots, \mathbf{x}_N) := \sum_{\substack{i=1 \\ i \neq j}}^{N} \sum_{j=1}^{N} \left[\mathrm{K}(\mathbf{x}_i, \mathbf{x}_j) + \mathrm{Q}(\mathbf{x}_i) + \mathrm{Q}(\mathbf{x}_j) \right]$$

among all sets of N points from Ω. The diagonal self-interaction terms are removed to allow singular kernels K. The external field Q is often taken to be zero. Evidently, further

requirements on the set Ω, the kernel K, and the external field Q are needed to ensure existence of a solution. A suitable Q compatible with K introduces soft boundaries and thus prevents points from escaping to infinity in the case of unbounded sets Ω. On the other hand, the restriction of the points to a finite set, fractal, torus or sphere of Hausdorff dimension $d \leq p$ introduces fractal or topological aspects. A standard assumption is that K is symmetric and lower semi-continuous on $\Omega \times \Omega$ and that Q is also lower semi-continuous on Ω. This implies that the minimal K-energy problem with external field Q has a solution for every $N \geq 2$ when solved for an infinite compact set Ω. A solution is called N-*point minimal* K-*energy configuration associated with* Q and its K-energy is equal to the *minimum* N-*point* K-*energy of* Ω *associated with* Q given by

$$\mathcal{E}_{\mathrm{K,Q}}(\Omega; N) := \inf \left\{ E_{\mathrm{K,Q}}(\mathbf{x}_1, \ldots, \mathbf{x}_N) : \mathbf{x}_1, \ldots, \mathbf{x}_N \in \Omega \right\}.$$

Furthermore, if $\mathcal{E}_{\mathrm{K,Q}}(\Omega; N) > 0$ for all N, then the quantities $N(N-1)/\mathcal{E}_{\mathrm{K,Q}}(\Omega; N)$ form a non-increasing sequence that is bounded from below by lower semi-continuity.[4] Thus the limit exists and it discriminates between two types of sets Ω depending on whether or not this limit vanishes. This gives rise to two different regimes characterised by a complete change in the nature of the minimisation problem with regard to properties of the solution and methods that are used to study it. In the field-free setting ($Q \equiv 0$), the normalised *discrete minimal* N-*point energy of* Ω, given by $\mathcal{E}_{\mathrm{K,0}}(\Omega; N)/[N(N-1)]$, is also known as the Nth diameter of Ω. The limit in the extended sense (as $N \to \infty$) is called the *transfinite diameter of* Ω.

A fundamental question concerns the "limit distribution" (if such exists) of a sequence (X_N^*) of minimal energy configurations X_N^* on Ω as $N \to \infty$; *i.e.*, is there a (unique) Borel probability measure μ_Ω^* supported on Ω that is the weak limit of the sequence formed by the discrete equal weight distribution supported on X_N^*,

$$\mu_{X_N^*} := \frac{1}{N} \sum_{j=1}^{N} \delta_{\mathbf{x}_j}.$$

Let $\mathcal{M}(\Omega)$ denote the collection of Borel probability measures supported on Ω. The analogous continuous energy problem is to find a measure in $\mathcal{M}(\Omega)$ that minimises the *weighted* K-*energy associated with* Q given by

$$\mathcal{I}_{\mathrm{K,Q}}(\mu) := \mathcal{I}_{\mathrm{K}}(\mu) + 2 \int_\Omega Q(\mathbf{x}) \, \mathrm{d}\mu(\mathbf{x})$$

among all measures in $\mathcal{M}(\Omega)$, where

$$\mathcal{I}_{\mathrm{K}}(\mu) := \int_\Omega \int_\Omega \mathrm{K}(\mathbf{x}, \mathbf{y}) \, \mathrm{d}\mu(\mathbf{x}) \mathrm{d}\mu(\mathbf{y})$$

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[4]The reciprocals $\mathcal{E}_{\mathrm{K,Q}}(\Omega; N)/[N(N-1)]$ always form a non-decreasing sequence that may be unbounded.

is the K-*energy of* $\mu \in \mathcal{M}(\Omega)$. A minimising measure $\mu_{K,Q;\Omega} \in \mathcal{M}(\Omega)$ with

$$\mathcal{I}_{K,Q}(\mu_{K,Q;\Omega}) = W_{K,Q}(\Omega) := \inf \left\{ \mathcal{I}_{K,Q}(\mu) \mid \mu \in \mathcal{M}(\Omega) \right\},$$

is called a K-*extremal (or positive equilibrium) measure on* Ω *associated with* Q. In the field free setting $Q \equiv 0$, a minimising measure $\mu_{K;\Omega}$ is called K-*equilibrium measure on* Ω. In this case, the K-energy $\mathcal{I}_K(\mu_{K;\Omega})$ is equal to the *Wiener energy of* Ω,

$$W_K(\Omega) := \inf \left\{ \mathcal{I}_K(\mu) \mid \mu \in \mathcal{M}(\Omega) \right\}.$$

A fundamental question concerns the relationship between the Wiener energy of Ω, the transfinite diameter of Ω, and the Chebyshev constant of Ω which is the limit as $N \to \infty$ of the Nth Chebyshev constant of Ω defined as

$$M_K(\Omega; N) := \sup_{\mathbf{x}_1, \ldots, \mathbf{x}_N \in \Omega} \inf_{\mathbf{x} \in \Omega} \frac{1}{N} \sum_{j=1}^{N} K(\mathbf{x}, \mathbf{x}_j);$$

see [93] and references cited therein for further details.

　　We conclude this section by a discussion of applications that make use of the minimal energy problem or are related to it.

The Thomson problem and its generalisations. A classical problem in electrostatics is to find the distribution of N unit point charges on a conductor in the most stable equilibrium (the charges interact according to the Coulomb potential $1/r$, where r is the Euclidean distance between two interacting charges). This leads to a minimisation problem for the potential energy of the discrete charge system named after J. J. Thomson who posed it for the sphere [194].[5] Generalisations of the Thomson problem that utilise a Riesz s-potential $1/r^s$ are used, *e.g.*, to model multi-electron bubbles and arrangements of protein subunits which form the shells (capsids) of viruses; see [38] (also [40, 41]) for a discussion. In material physics (see [63]) minimal energy points on the sphere have been used to model "spacer particles" in powders which ensure that spherical host-particles do not touch.

Polarisation. A problem related to finding the minimal K-energy configurations on a compact set $\Omega \subset \mathbb{R}^p$ is to find optimal N-*point K-polarisation configurations on* Ω, which are configurations on Ω that maximise the minimal value over Ω of the potential

$$\frac{1}{N} \sum_{j=1}^{N} K(\mathbf{x}, \mathbf{x}_j).$$

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[5]Recently, T. LaFave Jr. [141] investigated correspondences between the Thomson problem and atomic electronic structure, and he applied discrete transformations to the Thomson problem in [142] to study the minimal Coulomb energy.

An optimal configuration realises the Nth Chebyshev constant of Ω. For a Riesz kernel $K_s(\mathbf{x}, \mathbf{y}) = 1/[m(\mathbf{x}, \mathbf{y})]^s$ with metric m and $s > 0$, one has the following duality: the minimal s-energy configurations tend to *best-packing configurations* on Ω as $s \to \infty$, whereas the optimal s-polarisation configurations on Ω tend to *best-covering configurations* on Ω. Questions concerning optimal polarisation configurations, their limit distributions, polarisation inequalities, asymptotic behaviour of the Nth Chebyshev constant are addressed in [5, 32, 92, 117, 169]. The paper [92] proposes a conjecture for the dual of the Poppy-seed Bagel theorem[6] for optimal polarisation points (which is proven for the boundary case $s = \dim(\Omega)$ in [32]).

Smale's 7th Problem. In the seminal work [186], M. Shub and S. Smale define a condition number of a polynomial at a point in $\mathbb{C}$, which connects the problem of solving a polynomial equation with the discrete minimal logarithmic energy problem on $\mathbb{S}^2$ (see Section 3.2). They show that a monic polynomial whose zeros are the stereographic projection of minimal logarithmic energy points on the *Riemann sphere*, called "elliptic Fekete polynomial", is "well conditioned" and thus gives a good starting polynomial for a homotopy method for solving a polynomial equation or system (see [185]). This is the background for *Smale's 7th problem* [189]: *Find an algorithm which, on input N, outputs distinct points $\mathbf{x}_1, \ldots, \mathbf{x}_N$ on the Riemann sphere such that for a universal constant c,*

$$\sum_{\substack{i=1 \\ i \neq j}}^{N} \sum_{j=1}^{N} \log \frac{1}{\|\mathbf{x}_i - \mathbf{x}_j\|} - V_N \leq c \log N. \tag{3.1}$$

Here, V_N is the minimal logarithmic energy of N points on the Riemann sphere. Smale further specifies that the algorithm is a real number algorithm in the sense of Blum, Cucker, Shub, and Smale (see [27]) with halting time polynomial in N. In [19], C. Beltrán showed that there are N-point sets with logarithmic energy that differs from V_N by at most $1/9$. These points are rational with coordinates of order $\log_2 N$ bit length. Thus, there exists an exponential running time algorithm. Mean value considerations yield that the *typical* logarithmic energy of N i.i.d. uniformly distributed random points on the Riemann sphere is of the form $\frac{1}{2}N^2 - \frac{1}{2}N$. One the other hand, D. Armentano, C. Beltrán, and M. Shub [9] [7] observe that the typical logarithmic energy of the zeros of certain *random polynomials*, $\frac{1}{2}N^2 - \frac{1}{2}N \log N - \frac{1}{2}N$, is surprisingly small in the sense that the first two terms in the asymptotics of the minimal N-point logarithmic energy for the Riemann sphere are recovered. A recent account of the state of the art regarding Smale's 7th problem can be found in [20] (see also [21]).

Log gases, Coulomb gases, and random matrices. A log or Coulomb gas is a system

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[6]Cf. `http://news.vanderbilt.edu/2004/11/the-poppy-seed-bagel-theorem-59497/`
[7]The paper [97] gives a generalisation to higher dimensions and other manifolds.

of interacting particles in which the repelling interaction is governed by a logarithmic or Coulomb potential. An external field confines the particles to a finite volume of the space. Typically, the *mean-field regime* is considered. In this setting the number N of particles is large and the pair-interaction strength (coupling parameter) scales as the inverse of N. The Hamiltonian

$$H(\mathbf{x}_1, \ldots, \mathbf{x}_N) := \sum_{\substack{i=1 \\ i \neq j}}^{N} \sum_{j=1}^{N} k(\mathbf{x}_i - \mathbf{x}_j) + N \sum_{i=1}^{N} V(\mathbf{x}_i), \qquad \mathbf{x}_1, \ldots, \mathbf{x}_N \in \mathbb{R}^d,$$

where $k(\mathbf{x}) = -\log \|\mathbf{x}\|$ if $d = 2$ or $k(\mathbf{x}) = \|\mathbf{x}\|^{2-d}$ if $d \geq 3$, is minimised.[8] Very recent progress provides a deep connection with the discrete minimum energy problem on the sphere (see discussion at the end of Section 3.6). We refer further to S. R. Nodari and S. Serfaty [164] and N. Rougerie and S. Serfaty [174]. The concept of renormalised energy can be also successfully applied to random matrices; see A. Borodin and S. Serfaty [37]. For the connection between log gases and random matrix applications and theory we refer to the book of P. Forrester [98] and, *e.g.*, T. Claeys, A. B. J. Kuijlaars and M. Vanlessen [71] and A. Mays [155]. A related problem is the discrete energy of periodic point sets in the Euclidean space; see D. P. Hardin, E. B. Saff, and B. Simanek [120].

Half-toning. Loosely speaking, half-toning is a way of creating an illusion of a grey-value image by appropriately distributing black dots. In [111], M. Gräf, D. Potts, and G. Steidl show how the process of half-toning can be seen as a numerical integration problem. Here, the worst-case error can be interpreted as an external field problem where the picture drives the external field which guides the interacting points. The aim is then to minimise the worst-case error; see also [193] and [108].

Maximising Determinants. Points that maximise a Vandermonde-like determinant are well-suited for interpolation and numerical integration. They are called Fekete points due to the paper [96] by M. Fekete. Given a compact set in the complex plane, Fekete points are, indeed, minimal logarithmic energy points. However, in higher dimensions, Fekete's optimisation problem is different from minimising the logarithmic energy. I. H. Sloan and R. S. Womersley [187] used the logarithm of the determinant of an interpolation matrix to calculate extremal systems which yield interpolatory cubature rules with positive weights on the sphere. J. Marzo and J. Ortega-Cerdá [154] established that Fekete (or extremal) points are asymptotically uniformly distributed.

Diffusion on a sphere with localised traps. As an application in cellular signal transport, D. Coombs, R. Straube, and M. Ward [79] calculate the principal eigenvalue

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[8]Instead of the harmonic potentials one can also consider Riesz s-potentials $\|\cdot\|^{-s}$. For $s \geq d$ one needs to adjust the coupling parameter to ensure comparability between the pair-interaction part and external field part.

for the Laplacian on the unit sphere in the presence of N traps on the surface of the sphere of asymptotically small radii. The positions of the traps are chosen to minimise the discrete logarithmic energy given in (3.3) below.

3.2. The discrete and continuous minimal logarithmic and Riesz energy problem

The discrete logarithmic energy problem on $\mathbb{S}^d$ is concerned with the properties of N-point configurations $\{\mathbf{x}_1^*, \ldots, \mathbf{x}_N^*\} \subset \mathbb{S}^d$ that maximise the product of all mutual pairwise Euclidean distances

$$\prod_{\substack{i=1 \\ i \neq j}}^{N} \prod_{j=1}^{N} \|\mathbf{x}_i - \mathbf{x}_j\|, \tag{3.2}$$

or equivalently, minimise the discrete logarithmic energy

$$E_{\log}(\mathbf{x}_1, \ldots, \mathbf{x}_N) := \sum_{\substack{i=1 \\ i \neq j}}^{N} \sum_{j=1}^{N} \log \frac{1}{\|\mathbf{x}_i - \mathbf{x}_j\|} \tag{3.3}$$

over all N-point configurations $\{\mathbf{x}_1, \ldots, \mathbf{x}_N\}$ on $\mathbb{S}^d$. The discrete logarithmic energy can be understood as a limiting case (as $s \to 0$) of the Riesz s-energy

$$E_s(\mathbf{x}_1, \ldots, \mathbf{x}_N) := \sum_{\substack{i=1 \\ i \neq j}}^{N} \sum_{j=1}^{N} \frac{1}{\|\mathbf{x}_i - \mathbf{x}_j\|^s};$$

i.e., $E_s(\mathbf{x}_1, \ldots, \mathbf{x}_N) = N(N-1) + s\, E_{\log}(\mathbf{x}_1, \ldots, \mathbf{x}_N) + o(s)$ as $s \to 0$.

The discrete Riesz s-energy problem for $s > 0$ is concerned with the properties of N-point configurations $\{\mathbf{x}_1^*, \ldots, \mathbf{x}_N^*\} \subset \mathbb{S}^d$ that minimise the Riesz s-energy over all N-point configurations $\{\mathbf{x}_1, \ldots, \mathbf{x}_N\}$ on $\mathbb{S}^d$. For convenience we set $\mathrm{k}_s(\mathbf{x}, \mathbf{y}) := -\log\|\mathbf{x} - \mathbf{y}\|$ for $s = \log$ and $\mathrm{k}_s(\mathbf{x}, \mathbf{y}) := 1/\|\mathbf{x} - \mathbf{y}\|^s$ for $s \in \mathbb{R}$. Then we are interested in the *optimal N-point s-energy* of an infinite compact set $\Omega \subset \mathbb{S}^d$ defined by

$$\mathcal{E}_s(\Omega; N) := \begin{cases} \min\left\{ E_s(\mathbf{x}_1, \ldots, \mathbf{x}_N) \mid \mathbf{x}_1, \ldots, \mathbf{x}_N \in \Omega \right\} & \text{for } s = \log \text{ or } s \geq 0, \\ \max\left\{ E_s(\mathbf{x}_1, \ldots, \mathbf{x}_N) \mid \mathbf{x}_1, \ldots, \mathbf{x}_N \in \Omega \right\} & \text{for } s < 0. \end{cases}$$

Observe that $\mathcal{E}_0(\Omega; N) = N^2 - N$ (which is attained by any N-point set on Ω). The Riesz s-kernel is conditionally positive definite of order 1 for $-2 < s < 0$. Alternatively, as in the setting of numerical integration on $\mathbb{S}^d$, one can minimise

$$2 \int_\Omega \int_\Omega \|\mathbf{x} - \mathbf{y}\|^{-s}\, \mathrm{d}\sigma_d(\mathbf{x}) \mathrm{d}\sigma_d(\mathbf{y}) - \sum_{i=1}^{N} \sum_{j=1}^{N} \|\mathbf{x}_i - \mathbf{x}_j\|^{-s}$$

in this case. The papers [118, 176] are standard references for the discrete logarithmic and Riesz s-energy problem.

The *s-potential* and the *s-energy* of a measure μ in the class $\mathcal{M}(\Omega)$ of Borel probability measures supported on Ω are given, respectively, by

$$U_s^{\mu}(\mathbf{x}) := \int_{\Omega} \mathrm{k}_s(\mathbf{x}, \mathbf{y}) \, \mathrm{d}\mu(\mathbf{y}), \quad \mathbf{x} \in \mathbb{R}^{d+1}, \qquad \mathcal{I}_s(\mu) := \int_{\Omega} \int_{\Omega} \mathrm{k}_s(\mathbf{x}, \mathbf{y}) \, \mathrm{d}\mu(\mathbf{x}) \mathrm{d}\mu(\mathbf{y}). \quad (3.4)$$

For $s > 0$, the *s-capacity* of Ω is the reciprocal of the Wiener energy $\inf\{\mathcal{I}_s(\mu) \mid \mu \in \mathcal{M}(\Omega)\}$. When the Wiener energy is finite, it will be denoted by $W_s(\Omega)$. Because of the possibility of negative logarithmic energy integral, we set $\mathrm{cap}_{\log}(\Omega) := \exp(-\inf\{\mathcal{I}_{\log}(\mu) \mid \mu \in \mathcal{M}(\Omega)\})$. When finite, the infimum is denoted by $W_{\log}(\Omega)$. The lower semi-continuous logarithmic kernel is bounded from below and thus the kernel k_s (plus a constant) is strictly positive definite for $s = \log$ and $0 < s < d$. Consequently, the *s-equilibrium measure* $\mu_{\Omega,s}$ on Ω is unique for every compact set $\Omega \subset \mathbb{R}^p$ with finite *s*-energy; see [36, 143]. For the range $-2 < s < 0$, one also has a unique *s*-equilibrium measure on Ω; see [26] for the potential theoretic quantities and variational inequalities.

In the remaining section we consider the sphere $\mathbb{S}^d$ and subsets of $\mathbb{S}^d$. An *external field* is a lower semi-continuous function $\mathrm{Q} : \mathbb{S}^d \to (-\infty, \infty]$ such that $\mathrm{Q}(\mathbf{x}) < \infty$ on a set of positive Lebesgue surface measure. We note that the lower semi-continuity implies the existence of a finite c_{Q} such that $\mathrm{Q}(\mathbf{x}) \geq c_{\mathrm{Q}}$ for all $\mathbf{x} \in \mathbb{S}^d$. The *weighted energy associated with* $\mathrm{Q}(\mathbf{x})$ is then given by

$$\mathcal{I}_{\mathrm{Q},s}(\mu) := \mathcal{I}_s(\mu) + 2 \int \mathrm{Q}(\mathbf{x}) \mathrm{d}\mu(\mathbf{x}), \qquad \mu \in \mathcal{M}(\mathbb{S}^d). \quad (3.5)$$

We recall from [90] the following Frostman-type result that deals with existence and uniqueness of the *s*-extremal measure on Ω associated with Q and its characterisation in terms of weighted potentials. The result is stated for the Riesz case but a similar result holds also for the logarithmic case. The potential theory used in the context of this survey is formulated by G. Björck [26] (dealing with Riesz potential with negative exponent), by E. B. Saff, and V. Totik [177] (logarithmic external field problem in the plane), by N. S. Landkof [143] (Riesz and logarithmic potential and general reference) and, in particular, for Riesz external field problems by N. V. Zoriǐ [203, 204].

Proposition 3.1. *Let $0 < s < d$. For the minimal energy problem on $\mathbb{S}^d$ with external field Q the following properties hold:*

(a) $W_{\mathrm{Q},s} := \inf \{\mathcal{I}_{\mathrm{Q},s}(\mu) \big| \mu \in \mathcal{M}(\mathbb{S}^d)\}$ *is finite.*

(b) *There is a unique s-extremal measure $\mu_{\mathrm{Q},s} \in \mathcal{M}(\mathbb{S}^d)$ associated with Q. Moreover, the support $S_{\mathrm{Q},s} := \mathrm{supp}(\mu_{\mathrm{Q},s})$ of this measure is contained in the compact set $E_M := \{\mathbf{x} \in \mathbb{S}^d : \mathrm{Q}(\mathbf{x}) \leq M\}$ for some $M > 0$.*

(c) *The measure $\mu_{\mathrm{Q},s}$ satisfies the variational inequalities*

$$U_s^{\mu_{\mathrm{Q},s}}(\mathbf{x}) + \mathrm{Q}(\mathbf{x}) \geq F_{\mathrm{Q},s} \quad q.e. \ on \ \mathbb{S}^d, \quad (3.6)$$
$$U_s^{\mu_{\mathrm{Q},s}}(\mathbf{x}) + \mathrm{Q}(\mathbf{x}) \leq F_{\mathrm{Q},s} \quad everywhere \ on \ S_{\mathrm{Q},s}, \quad (3.7)$$

where

$$F_{Q,s} := W_{Q,s} - \int Q(\mathbf{x}) \mathrm{d}\mu_{Q,s}(\mathbf{x}). \tag{3.8}$$

(d) *Inequalities* (3.6) *and* (3.7) *completely characterise the s-extremal measure* μ_Q *in the sense that if* $\nu \in \mathcal{M}(\mathbb{S}^d)$ *is a measure with finite s-energy such that for some constant* C *we have*

$$U_s^\nu(\mathbf{x}) + Q(\mathbf{x}) \geq C \quad q.e. \ on \ \mathbb{S}^d, \tag{3.9}$$
$$U_s^\nu(\mathbf{x}) + Q(\mathbf{x}) \leq C \quad everywhere \ on \ \mathrm{supp}(\nu), \tag{3.10}$$

then $\nu = \mu_{Q,s}$ *and* $C = F_{Q,s}$.

A property holds *quasi-everywhere* if the exceptional set has s-capacity zero. If Q is continuous on $\mathbb{S}^d$, then the inequalities (3.6) and (3.9) hold everywhere on $\mathbb{S}^d$.

In principle, once $\mathrm{supp}(\mu_Q)$ is known, then the measure μ_Q can be recovered by solving an integral equation for the weighted s-potential arising from (3.9) and (3.10). Finding $\mathrm{supp}(\mu_Q)$ when it is a proper subset of $\mathbb{S}^d$ can be a very difficult problem. It is a substantially easier task to find a signed measure that has constant weighted s-potential everywhere on $\mathbb{S}^d$. Given a compact subset $\Omega \subset \mathbb{S}^d$ with $\mathrm{cap}_s(\Omega) > 0$ and an external field Q, there is a unique (if it exists, see [52, Lemma 23]) *signed s-equilibrium* measure $\eta_{K,Q}$ supported on Ω of total charge one associated with Q with constant weighted s-potential; *i.e.*,

$$U_s^{\eta_{K,Q}}(\mathbf{x}) + Q(\mathbf{x}) = G_{K,Q,s} \qquad \text{everywhere on } \Omega$$

for some constant $G_{K,Q,s}$. A remarkable connection exists to the analogue of the *Mhaskar-Saff functional* from classical planar potential theory ([157] and [177, Chapter IV, p. 194]) given by

$$\mathcal{F}_s(\Omega') := W_s(\Omega') + \int Q(\mathbf{x}) \, \mathrm{d}\mu_{\Omega'}(\mathbf{x}), \qquad \Omega' \subset \mathbb{S}^d \text{ compact with } \mathrm{cap}_s(\Omega') > 0,$$

where $W_s(\Omega')$ is the s-energy of Ω' and $\mu_{\Omega'}$ is the s-equilibrium measure (without external field) on Ω'. Namely, if the signed s-equilibrium on a compact set Ω' associated with Q exists, then $\mathcal{F}_s(\Omega') = G_{K,Q,s}$. The essential property of the $\mathcal{F}_s$-functional is that it is minimised for the support of the s-extremal measure (see [53, Proposition 8] for a precise statement). The papers [52, 53, 54, 90] determine the signed s-equilibrium on the full sphere and on spherical caps associated with logarithmic and Riesz s-external fields due to a single point charge (or a axis-supported superposition of such fields). This signed s-equilibrium characterises then the s-extremal measure on $\mathbb{S}^d$. A further application is the proof of optimal separation ("well-separation") for minimal Riesz s-energy points on $\mathbb{S}^d$ for the range $s \in [d-2, d)$.[9] The use of balayage techniques (the signed equilibrium can be

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[9]For further discussion of the separation of minimal s-energy points on $\mathbb{S}^d$, see [52] and citations therein. It should be mentioned that S. B. Damelin and V. Maymeskul [86] give a separation result of order $N^{-1/(s+2)}$, $0 < s \leq d-2$, which is sharp for $s = d-2 > 0$.

expressed as the difference of two balayage measures) together with a restricted principle of domination and a restricted maximum principle yields that the Riesz parameter s is restricted to the interval $[d-2, d)$. New phenomena occur when $s = d - 2$. For example, a boundary charge appears in the signed $(d-2)$-equilibrium for a spherical cap which vanishes when the cap coincides with the support of the s-extremal measure on $\mathbb{S}^d$. Numerical methods for external field problems for Riesz potentials have been developed and applied in [114, 115, 116, 166].

3.3. The distribution of minimal logarithmic and Riesz energy points

Let Ω be an infinite compact set $\Omega \subset \mathbb{R}^p$ with Hausdorff dimension d having finite logarithmic or Riesz s-energy $(s > 0)$ and $Q \equiv 0$. Then classical potential theory implies that the minimal energy configurations X_N^* on Ω are distributed according to the unique equilibrium measure μ_Ω on Ω and the discrete measures $\mu_{X_N^*}$ have μ_Ω as a weak limit. The s-equilibrium measure on $\mathbb{S}^d$ is the uniform measure σ_d. In general, the measure μ_Ω will not be uniform as the example of a circular torus shows (see [121] and [55, 56]).[10] In this potential theoretic regime $(s = \log \text{ or } 0 < s < d)$, the support of μ_Ω depends on whether the kernel is superharmonic, harmonic or subharmonic (see [143]). In the superharmonic case $(s = \log \text{ or } 0 < s < p - 2)$, the measure μ_Ω is supported on the *outer boundary* (*i.e.*, the boundary of Ω shared with the unbounded component of the ambient space $\mathbb{R}^p$), whereas in the strictly subharmonic case, the measure μ_Ω can be supported on all of Ω.

The intuition is that in the regime $s = \log \text{ or } 0 < s < d$ global effects dominate (points interact as if they are subject to long-range forces and the range increases as s becomes smaller). In the hypersingular case $s > d$, local effects dominate (points interact as if they are responding to a short range force). Both kinds of interactions intermingle when $s = d$.

In the hypersingular case $s \geq d$, the energy integral attains $+\infty$ for every $\mu \in \mathcal{M}(\Omega)$. Geometric measure theory yields that the limiting distribution of a sequence (X_N^*) of minimal Riesz s-energy N-point sets on Ω (even asymptotically s-energy minimising would suffice) is uniformly distributed with respect to the d-dimensional Hausdorff measure $\mathcal{H}_d$,

$$\mu_{X_N^*} \overset{*}{\longrightarrow} \frac{\mathcal{H}_d\big|_\Omega}{\mathcal{H}_d(\Omega)} \qquad \text{as } N \to \infty,$$

for a large class of sets Ω with $\mathcal{H}_d(\Omega) > 0$ (see [34] and earlier work [33, 118, 119]). It should be noted that in the case $s = d$ an additional regularity assumption on Ω is required. In the limit $s \to \infty$ only the nearest-neighbour interaction matters and the optimal solutions are best-packing configurations which solve the Tammes problem [192]. The paper [34] also shows analogous results for weighted Riesz s-energy

$$\sum_{\substack{i=1 \\ i \neq j}}^{N} \sum_{j=1}^{N} \frac{w(\mathbf{x}_i, \mathbf{x}_j)}{\|\mathbf{x}_i - \mathbf{x}_j\|^s},$$

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[10]It seems to be unresolved for which Ω the s-equilibrium measure on Ω is the uniform measure on Ω.

where w is (almost everywhere) **c**ontinuous and **p**ositive on the **d**iagonal (called CPD weight function). We remark that the case $s = d$ for $\mathbb{S}^d$ has already been dealt with in [103] using results from [139]. Furthermore, in [66] it was shown that the s-equilibrium measures on Ω converge to the normalised d-dimensional Hausdorff measure restricted to Ω as $s \to d^-$ under rather general assumptions on Ω. We remark that I. Pritzker [168] studied the discrete approximation of the equilibrium measure on a compact set $\Omega \subset \mathbb{R}^p$, $p \geq 2$, with positive s-capacity by means of points which do not need to lie inside Ω. He also obtained discrepancy estimates in the harmonic case. The properties of so-called *greedy* K-*energy points* were studied by A. López García and E. B. Saff [148].

Summarising, for $-2 < s < 0$, $s = \log$ and $s > 0$, the (asymptotically) s-energy minimising N-point configuration on $\mathbb{S}^d$ are uniformly distributed with respect to the surface area measure σ_d.

In certain applications one prefers to generate well distributed N-point sets on a compact d-rectifiable set in $\mathbb{R}^p$ which have a prescribed non-uniform asymptotic distribution ρ with respect to $\mathcal{H}_d$ as $N \to \infty$. It is shown in [122] that such points can be obtained by minimising the energy of N points on Ω interacting via a weighted power law potential $w(\mathbf{x}, \mathbf{y})/\|\mathbf{x} - \mathbf{y}\|^s$, where $s > d$ and $w(\mathbf{x}, \mathbf{y}) := [\rho(\mathbf{x})\rho(\mathbf{y})]^{-s/(2d)}$. Furthermore, such point sets are "quasi-uniform" in the sense that the ratio of the covering radius to the separation distance is uniformly bounded in N.[11] As mentioned in the introduction, quasi-uniformity is crucial for a number of numerical methods (see [99, 144, 180]). S. V. Borodachov, D. P. Hardin, and E. B. Saff show in [35] that it suffices to use a varying truncated weight $w(\mathbf{x}, \mathbf{y}) \Phi(\|\mathbf{x} - \mathbf{y}\|/r_N)$. Thus only those pairs of points that are located at a distance of at most $r_N = C_N N^{-1/d}$ from each other contribute to the energy sum. (The positive sequence (C_N) can be taken to tend to ∞ as slowly as one wishes.) In this way, under suitable assumptions, the complexity of the energy computation can be greatly reduced, leading to order $N C_N^d$ computations for generating "low energy" N-point approximations.[12]

A point charge approaching the sphere subject to the same law of interaction as the points on the sphere affects the charge distribution on the sphere. Sufficiently close to the sphere, it will generate a spherical cap with zero charge. The papers [52, 53, 54], in particular, provide explicit representations of the charge distributions due to a single external charge. They also address a question attributed to A. A. Gonchar, namely to find a critical distance from the sphere surface of a point charge generating the external field so that the support of the s-extremal measure on $\mathbb{S}^d$ only just becomes the whole sphere. In the harmonic case this distance is characterised by the largest zero of polynomials dubbed Gonchar polynomials.[13]

For a small N, numerical optimisation methods can be used to find putative minimal

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[11]For the constant weight case, see also [86].

[12]We note that in the astronomical community a hierarchical equal area iso-latitude pixelisation (HEALPix [100]) is used to generate large numbers of uniformly distributed points on $\mathbb{S}^2$; cf. also [145]. A. Holhoş and D. Roşca [133] study Riesz energy of points derived from an area-preserving map from the 2-sphere to the octahedron.

[13]For $d = 2$ and $d = 4$, the critical distance is the golden ratio and the plastic number.

Riesz s-energy configurations. T. Erber and G. M. Hockney [91] noticed that the number of local minimal energy configurations (most of which are not global ones) seems to grow exponentially with N. Most of the numerical data is for the 2-sphere and for the Coulomb case and a few other values of s. Regarding data, we refer to online resources [39, 127, 201] and a more recent study of the energy landscape in [64, 65], whereas [22] provides a complexity analysis for the logarithmic case. Higher dimensional configurations have also been investigated (see [13]). In [163] monotonicity properties of the second discrete derivative were considered and led to new putative low-energy configurations in two cases. Numerical results in [156] suggest that a minimal s-energy N-point set may transit through one or more (depending on N) basic configuration as s grows.[14] The smallest only partly resolved problem is the *five point problem on $\mathbb{S}^2$*. Five points cannot form a universally optimal system [74].[15] Melnyk *et al.* [156] identified two basic configurations: triangular bi-pyramid and quadratic pyramid. According to numerical results, the regular triangular bi-pyramid is the putative energy-minimising configuration for $2 \leq s \leq 15.048077392\ldots$, whereas for higher values of s it seems to be the square pyramid (with adjusted height); see also [163] for a finer analysis. Moreover, it is shown in [28] that there are sequences of s-energy minimising configurations that tend to a square pyramid best packing configuration as $s \to \infty$. In general, the five point problem is a difficult problem to analyse rigorously. Recently, the papers [182] (for the Coulomb case $s = 1$ and for $s = 2$) and [134] (for sum of distances, $s = -1$) provided computer-assisted proofs that the triangular bi-pyramid is optimal, whereas in the logarithmic case a conventional proof was given in [89]. In [197] a bi-quadratic energy functional is considered. Other rigorously proved minimising configurations are rare and are often universally optimal. Proved minimising configurations on $\mathbb{S}^2$ are the antipodal and equilateral configuration and the Platonic solids with $N = 4, 6$ and 12 vertices. For higher dimensions, we refer to [74, Table 1], [11] and [77, Section 5.3].

3.4. Asymptotic expansion of minimal Riesz energy

Let $s > 0$. As sets we shall consider the unit sphere $\mathbb{S}^d$ and, more generally, infinite compact sets $\Omega \subset \mathbb{R}^p$ with Hausdorff dimension $0 < d \leq p$. The leading term of the asymptotic expansion of the N-point Riesz s-energy of Ω is well-understood if Ω has positive s-capacity (*i.e.*, finite Riesz s-energy). This is the *potential-theoretic regime*. A standard argument from classical potential theory yields that the positive quantities $N(N - 1)/\mathcal{E}_s(\Omega; N)$ form a monotonically decreasing sequence. The limit $\text{diam}_s(\Omega)$, called the *generalised transfinite diameter of* Ω, is equal to the s-capacity of Ω (cf. [167]). Thus, the leading term of $\mathcal{E}_s(\Omega; N)$ grows like N^2 as $N \to \infty$ and the leading coefficient is given by the Riesz s-energy of Ω,

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[14]The phenomenon of transiting through several basic types of configurations can also be observed in the external field setting when s is fixed but the distance of the external field source varies (see [53, Figure 4]).

[15]A *universally optimal* point set minimises all energy functionals with kernels of the form $K(\mathbf{x}, \mathbf{y}) = f(\|\mathbf{x} - \mathbf{y}\|^2)$, where f is a completely monotonic C^∞ function like the Riesz potential $1/r^s$ for $s > 0$, meaning that $(-1)^k f^{(k)}(x) \geq 0$ for all k; see [74].

or equivalently, by the reciprocals of the s-capacity and transfinite diameter of Ω:

$$\lim_{N \to \infty} \frac{\mathcal{E}_s(\Omega; N)}{N^2} = W_s(\Omega) = \frac{1}{\mathrm{cap}_s(\Omega)} = \frac{1}{\mathrm{diam}_s(\Omega)}. \tag{3.11}$$

For $0 < s < d$, the Riesz s-energy of the sphere $\mathbb{S}^d$ has the explicit form

$$W_s(\mathbb{S}^d) = \mathcal{I}_s[\sigma_d] = 2^{d-1-s} \frac{\Gamma((d+1)/2)\,\Gamma((d-s)/2)}{\sqrt{\pi}\,\Gamma(d-s/2)}, \tag{3.12}$$

expressed in terms of the gamma function Γ. By identifying $W_s(\mathbb{S}^d)$ with the analytic continuation of the right-hand side above to the complex s-plane[16], we can define the Riesz s-energy of $\mathbb{S}^d$ for Riesz parameter s for which the s-energy integral (3.4) is $+\infty$ for every Borel probability measure on $\mathbb{S}^d$. The combined effort of [43, 139, 170, 199, 200] resulted in the following bounds for the second term of the minimal energy asymptotics:[17] there exist constants $c, C > 0$ depending only on $d \geq 2$ and $0 < s < d$ such that

$$c\,N^{1+s/d} \leq \mathcal{E}_s(\mathbb{S}^d; N) - W_s(\mathbb{S}^d)\,N^2 \leq C\,N^{1+s/d}, \qquad N \geq 2.$$

These estimates give the correct order of growth and sign for the second-order term. It is an open problem if the sequence $\left(\mathcal{E}_s(\mathbb{S}^d; N) - W_s(\mathbb{S}^d)\,N^2\right)/N^{1+s/d}$ has a limit as $N \to \infty$. A. A. Berezin [24] used a semi-continuum approach (a classical method from solid state physics, cf. [104]) to derive the plausible asymptotics

$$\mathcal{E}_s(\mathbb{S}^2; N) \approx N^2 \frac{2^{1-s}}{2-s} \left[1 - (n/N)^{1-s/2} \right] + N^{1+s/2} \left(\frac{\sqrt{3}}{8\pi} \right)^{s/2}$$

$$\times \left\{ \frac{6}{1^s} + \frac{6}{(\sqrt{3})^s} + \frac{6}{2^s} + \frac{12}{(\sqrt{7})^s} + \frac{6}{3^s} + \frac{6}{(2\sqrt{3})^s} + \frac{12}{(\sqrt{13})^s} + \cdots \right\}$$

based on the assumptions that a typical point (and most of its immediate neighbours) in a minimal Riesz s-energy N-point configuration on $\mathbb{S}^2$ gives rise to a hexagonal Voronoi cell (sixfold symmetry) whereas the defects according to the curved surface of the sphere will have no significant influence on the second term in the asymptotics. Thus the contribution to the Riesz s-energy due to a typical point can be split into a local part which uses n nearest neighbour points from a suitably adjusted flat hexagonal lattice and a distant part where the $N - n$ points are replaced by the continuous uniform distribution. The expression in curly braces gives the formal series expansion of the Epstein zeta function of the hexagonal lattices truncated to include only the $n - 1$ shortest distances in the lattice[18]. (In order to get a non-trivial expansion, n has to grow slowly in terms of N to

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[16]The meromorphic function $W_s(\mathbb{S}^d)$, which appears in the conjecture for the asymptotics in the hypersingular case, has simple poles (finitely many if d is even and infinitely many if d is odd). The effect of this dichotomy on the asymptotic expansion of the minimal Riesz s-energy is completely open for $d \geq 2$ and leads to $\log N$ terms for the unit circle, cf. [57].

[17]Similar estimates but with negative constants c, C hold for the sum of generalised distances (*i.e.*, $-2 < s < 0$); see [2, 3, 4, 18, 128, 190, 191] and culminating in [199, 200].

[18]Indeed, the first few most frequent distances in a putative minimal energy configuration emulate remarkably well the first few distances in a hexagonal lattice (see, in particular, [59, Figure 1]).

infinity. The paper [24] mentions some numerical experiments for slowly growing n but a rigorous investigation has not been undertaken.) The discussion leading to Conjecture 1 below suggests that the semi-continuum approach would also work for $d = 4$, 8, and 24. In general, it is not clear which local approximation should be used.

For Ω in $\mathbb{R}^p$ with vanishing s-capacity, the leading term is rather well-understood. In the *strictly hypersingular regime* $s > d$, Hardin and Saff [119] (for rectifiable d-dimensional manifolds including the sphere $\mathbb{S}^d$) and Borodachov, Hardin, and Saff [34] (for infinite compact d-rectifiable sets[19]) established the existence of a constant $C_{s,d}$ such that for a large class of sets Ω [20]

$$\lim_{N \to \infty} \frac{\mathcal{E}_s(\Omega; N)}{N^{1+s/d}} = \frac{C_{s,d}}{[\mathcal{H}_d(\Omega)]^{s/d}}, \tag{3.13}$$

where $\mathcal{H}_d$ denotes the d-dimensional Hausdorff measure in $\mathbb{R}^p$ normalised such that the d-dimensional unit cube has $\mathcal{H}_d$-measure 1. This result is referred to as the *Poppy-seed Bagel Theorem* because of its interpretation for distributing points on a torus. Except for one-dimensional sets (when $C_{s,1}$ is twice the Riemann zeta function at s, see [153, Thm. 3.1]), the precise value of $C_{s,d}$ is not known. Its determination is a challenging open problem. The significance and difficulty of obtaining $C_{s,d}$ is due to the deep connection to densest packings. In [33] it is shown that $C_{s,d}$ is tied to the largest sphere packing density Δ_d in $\mathbb{R}^d$ and the best-packing distance δ_N^* of N-points on $\mathbb{S}^d$ by means of the limit relations[21]

$$\lim_{s \to \infty} [C_{s,d}]^{-1/s} = 2 \left[\frac{\Delta_d}{\mathcal{H}_d(\mathbb{B}^d)} \right]^{1/d} = \lim_{N \to \infty} N^{1/d} \delta_N^*. \tag{3.14}$$

Here, $\mathcal{H}_d(\mathbb{B}^d)$ is the volume of the unit ball in $\mathbb{R}^d$. We recall that Δ_d is only known for three cases: $\Delta_1 = 1$, $\Delta_2 = \pi/\sqrt{12}$ (Thue [195] and L. Fejes Tóth [95]) and $\Delta_3 = \pi/\sqrt{18}$ (Kepler conjecture proved by Hales [113]). The connection to (regular) lattices comes from the fact that the Riesz s-energies of shrunk copies of a lattice Λ in $\mathbb{R}^d$ restricted to the fundamental parallelotope Ω of Λ yield an upper estimate for $C_{s,d}$ (for $s > d$): consider the $N = n^d$ points $X_N = \frac{1}{n}\Lambda \cap \Omega$, then one has (cf. [59, Prop. 1])

$$\mathcal{E}_s(\Omega; N) \leq E_s(X_N) \leq n^{d+s} \zeta_\Lambda(s) = N^{1+s/d} \zeta_\Lambda(s).$$

This implies

$$C_{s,d} \leq \min_\Lambda |\Lambda|^{s/d} \zeta_\Lambda(s), \tag{3.15}$$

where the minimum is extended over all lattices Λ in $\mathbb{R}^d$ with positive co-volume $|\Lambda|$. Because of (3.14), the sharpness of this inequality touches on questions regarding densest lattice sphere packings. For $1 \leq d \leq 8$ and $d = 24$, the unique densest lattice in $\mathbb{R}^d$ up to scaling and isometries is the root lattice A_1, A_2, A_3, D_4, D_5, E_6, E_7, E_8, and the Leech

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[19]A d-rectifiable set is the Lipschitz image of a bounded set in $\mathbb{R}^d$.

[20]The boundedness of $\mathcal{E}_s(\mathbb{S}^d; N)/N^{1+s/d}$ has already been shown in [139].

[21]Indeed, one can recast this relation as $\Delta_d = \lim_{s \to \infty} \lim_{N \to \infty} \left[\mathcal{E}_s(\frac{1}{2}\mathbb{B}^d; N)/N^{1+s/d} \right]^{-d/s}$.

lattice Λ_{24}, respectively (cf. [75]). Among those the hexagonal lattice A_2 in $\mathbb{R}^2$, the E_8 root lattice in $\mathbb{R}^8$ and the Leech lattice Λ_{24} in $\mathbb{R}^{24}$ are conjectured to be universally optimal, whereas the remaining lattices are provably not universally optimal (cf. [74, 76]). See [179] for local optimality results and [81] for improvements. Montgomery [158] proved that the hexagonal lattice is universally optimal amongst all lattices in $\mathbb{R}^2$ (which is weaker than universal optimality amongst all periodic point configurations). H. Cohn and N. Elkies [73] conjectured that E_8 and the Leech lattice Λ_{24} solve the sphere packing problem in their dimension. It is generally expected that for sufficiently large d, lattice packings are not densest packings and [196] suggests that best-packings are highly "disordered" as $d \to \infty$. This motivates the following conjecture[22].

Conjecture 1 ([57]). *For $d = 2$, 4, 8, and 24, one has $C_{s,d} = |\Lambda_d|^{s/d} \zeta_{\Lambda_d}(s)$, where Λ_d denotes, respectively, the hexagonal lattice A_2, the root lattices D_4 and E_8, and the Leech lattice Λ_{24}.*

We remark that in [59] very coarse lower and upper bounds are obtained for $\mathcal{E}_s(\mathbb{S}^d; N)$ which imply for $s > d \geq 2$ and $(s - d)/2$ not an integer, the estimates

$$\frac{d}{s-d} \left[\frac{1}{2} \frac{\Gamma((d+1)/2)\,\Gamma(1+(s-d)/2)}{\sqrt{\pi}\,\Gamma(1+s/2)} \right]^{s/d} \leq \frac{C_{s,d}}{[\mathcal{H}_d(\mathbb{S}^d)]^{s/d}} \leq \left[\frac{\mathcal{H}_d(\mathbb{B}^d)}{\mathcal{H}_d(\mathbb{S}^d)(1-d/s)} \right]^{s/d}.$$

In the *hypersingular case* $s = d$, more can be said. It has been known from [139] that the leading term of $\mathcal{E}_d(\mathbb{S}^d; N)$ grows like $N^2 \log N$ and

$$\lim_{N \to \infty} \frac{\mathcal{E}_d(\mathbb{S}^d; N)}{N^2 \log N} = \frac{\mathcal{H}_d(\mathbb{B}^d)}{\mathcal{H}_d(\mathbb{S}^d)} = \frac{1}{d} \frac{\omega_{d-1}}{\omega_d} = \frac{1}{d} \frac{\Gamma((d+1)/2)}{\sqrt{\pi}\,\Gamma(d/2)}. \tag{3.16}$$

The best estimate so far for the second-order term has been obtained recently in [59],

$$-c(d)\,N^2 + \mathcal{O}(N^{2-2/d}\log N) \leq \mathcal{E}_d(\mathbb{S}^d; N) - \frac{\mathcal{H}_d(\mathbb{B}^d)}{\mathcal{H}_d(\mathbb{S}^d)}\,N^2 \log N \leq \frac{\mathcal{H}_d(\mathbb{B}^d)}{\mathcal{H}_d(\mathbb{S}^d)}\,N^2 \log\log N + \mathcal{O}(N^2)$$

as $N \to \infty$, where the constant $c(d)$ is given by

$$c(2) = 1/2, \qquad c(d) := \frac{\mathcal{H}_d(\mathbb{B}^d)}{\mathcal{H}_d(\mathbb{S}^d)} \left\{ 1 - \log \frac{\mathcal{H}_d(\mathbb{B}^d)}{\mathcal{H}_d(\mathbb{S}^d)} + d\left[\psi(d/2) - \psi(1) - \log 2\right] \right\} > 0.$$

(Recall, that $\psi := \Gamma'/\Gamma$ denotes the digamma function.) Based on a limiting process $s \to d$ in Conjecture 2 given below, it is conjectured in [59] that the correct order of the second term is N^2. Furthermore, in the case $d = 2$ a conjecture is posed for the constant of the N^2-term.

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[22]The conjecture for $d = 2$ appeared in [139].

3.5. Higher Order Terms – Complete Asymptotic Expansions – Fundamental Conjecture

Very little is known about higher-order terms of the asymptotics of the minimal Riesz s-energy except for the unit circle. As the Nth roots of unity are universally optimal,[23] the complete asymptotic expansion can be obtained by direct computation of the Riesz s-energy $\mathcal{L}_s(N)$ of the Nth roots of unity, see [57] for Euclidean and [58] for geodesic metric. Indeed, for $s \in \mathbb{C}$ with $s \neq 0, 1, 3, 5, \ldots$ and fixed $p = 1, 2, 3, \ldots$, one has[24]

$$\mathcal{L}_s(N) = W_s(\mathbb{S}^1)\,N^2 + \frac{2\,\zeta(s)}{(2\pi)^s}\,N^{1+s} + \sum_{n=1}^{p} \alpha_n(s)\frac{2\,\zeta(s-2n)}{(2\pi)^s}\,N^{1+s-2n}$$
$$+ \mathcal{O}_{s,p}(N^{-1+\mathrm{Re}(s)-2p}) \qquad \text{as } N \to \infty, \tag{3.17}$$

where the coefficients $\alpha_n(s)$, $n \geq 0$, are defined by the generating function relation

$$\left(\frac{\sin \pi z}{\pi z}\right)^{-s} = \sum_{n=0}^{\infty} \alpha_n(s)\,z^{2n}, \quad |z| < 1, \ s \in \mathbb{C}.$$

Explicit formulas for $\alpha_n(s)$ in terms of generalised Bernoulli polynomials $B_n^{(\alpha)}(x)$ are given in [58]. The asymptotics (3.17) has two noteworthy features: this expansion is valid for complex s and the Riemann zeta function plays an essential role. The coefficients of the terms in the asymptotics are best understood as functions in the complex s-plane. This is called the *principle of analytic continuation.*[25] The interplay between the simple poles of the coefficient $W_s(\mathbb{S}^1)$ of the N^2-term in (3.12) and the simple poles of the shifted Riemann zeta functions then gives rise to a logarithmic term whenever s tends to one of the exceptional cases $s = 0, 1, 3, 5, \ldots$.

By combining the results for the potential theoretic and the hypersingular regime, the principle of analytic continuation motivates the following fundamental conjecture.

Conjecture 2 (see [59]). *Let $d \geq 2$. Then for $0 < s < d+2$ with $s \neq d$,*

$$\mathcal{E}_s(\mathbb{S}^d; N) = W_s(\mathbb{S}^d)\,N^2 + \frac{C_d(s)}{[\mathcal{H}_d(\mathbb{S}^d)]^{s/d}}\,N^{1+s/d} + o(N^{1+s/d}) \qquad \text{as } N \to \infty,$$

where $W_s(\mathbb{S}^d)$ is the analytic continuation of the right-hand side of (3.12) and $C_d(s)$ is the analytic continuation of $C_{s,d}$ in (3.13). Furthermore, for $d = 2, 4, 8$, and 24, the constant $C_d(s)$ is the analytic continuation of $|\Lambda_d|^{s/d}\,\zeta_{\Lambda_d}(s)$, where Λ_d is given in Conjecture 1.

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[23]For $s \geq -1$ (and $s \neq 0$) a convexity argument can be applied to get optimality for Riesz s-energy (see [4, 94, 102]). The much more general result [74, Theorem 1.2] provides optimality for $s > -2$.

[24]The precise formulas for finite $N \geq 2$ are obtained in [46].

[25]This principle breaks down when the perfectly symmetric unit circle is replaced by some other smooth closed curve Γ. Then the s-equilibrium measure on Γ is not the normalised arc-length measure for each $0 < s < 1$ which plays a role in the characterisation of the coefficient of N^2 in the hypersingular regime $1 < s < 3$; see [31].

It should be discussed briefly that the asymptotic expansion of the minimal s-energy can also be studied from a geometrical point of view by identifying which features of the Voronoi cell decomposition (or its dual, the Delaunay triangulation) induced by minimal s-energy configurations contribute in which way to the asymptotics. For the 2-sphere and large N, the typical picture is a vast sea of hexagonal Voronoi cells — thus the local approximation of the neighbourhood of a typical point is done by a suitably scaled hexagonal lattice to get the second term of the asymptotics. The topology of the sphere gives rise to *geometric frustration* (cf. [175]) where certain points pick up a *topological charge* that measures the discrepancy from the ideal coordination number (six) of the planar triangular lattice. Euler's celebrated Polyhedral formula yields that the total topological charge on $\mathbb{S}^2$ is always 12. Numerically, one observes "scars" for large N emerging from 12 pentagonal centres. These scars attract pentagon-heptagon pairs which have total topological charge zero. It is an unresolved question if there are Voronoi cells with more than 7 sides in minimising configurations. It is not well understood how (*i.e.*, on which level of the asymptotic scale) scars, the type of Voronoi cells, and the variation in their sizes affect higher-order terms of the asymptotics; see [65] for numerical background and [41] for an approach using elastic continuum formalism.

A much harder question concerns the (asymptotic) behaviour of the "point-energies" of an optimal configuration (*i.e.*, the contributions to the s-energy due to individual points of the configuration — the s-energy is the total sum of these point-energies). Point energies are considered in [65] (see Figure 1) and asymptotic estimates are given in [86].

3.6. Asymptotic expansion of logarithmic energy

The leading term of the asymptotic expansion for a compact set Ω in $\mathbb{R}^p$ with positive logarithmic capacity (*i.e.*, finite logarithmic energy) follows from classical potential theory. It should be noted that (see [36])

$$\frac{\mathrm{d}}{\mathrm{d}s}\mathcal{E}_s(\Omega; N)\bigg|_{s=0^+} = \mathcal{E}_{\log}(\Omega; N), \qquad N \geq 2. \qquad (3.18)$$

For the unit sphere $\mathbb{S}^d$, one has

$$\lim_{N\to\infty} \frac{\mathcal{E}_{\log}(\mathbb{S}^d; N)}{N^2} = W_{\log}(\mathbb{S}^d) = \log \frac{1}{\mathrm{cap}_{\log}(\mathbb{S}^d)},$$

where the *logarithmic energy of* $\mathbb{S}^d$ is given by

$$W_{\log}(\mathbb{S}^d) = \frac{\mathrm{d}W_s(\mathbb{S}^d)}{\mathrm{d}s}\bigg|_{s=0^+} = \log\frac{1}{2} + \frac{1}{2}\left[\psi(d) - \psi(d/2)\right].$$

An averaging argument that uses an equal-area partition of $\mathbb{S}^d$ and bounds of G. Wagner [200] and [44] yields

$$\mathcal{E}_{\log}(\mathbb{S}^d; N) = W_{\log}(\mathbb{S}^d)\, N^2 - \frac{1}{d}\, N \log N + \mathcal{O}(N), \qquad N \to \infty.$$

Relation (3.18) and Conjecture 2 provide the basis for the following conjecture posed in [59].

Conjecture 3. *For $d = 2$, 4, 8, and 24,*

$$\mathcal{E}_{\log}(\mathbb{S}^d; N) = W_{\log}(\mathbb{S}^d) \, N^2 - \frac{1}{d} N \log N + C_{\log,d} \, N + o(N) \qquad \textit{as } N \to \infty,$$

where

$$C_{\log,d} = \frac{1}{d} \log \frac{\mathcal{H}_d(\mathbb{S}^d)}{|\Lambda_d|} + \zeta'_{\Lambda_d}(0).$$

For $d = 2$ one has

$$C_{\log,2} = 2 \log 2 + \frac{1}{2} \log \frac{2}{3} + 3 \log \frac{\sqrt{\pi}}{\Gamma(1/3)} = -0.05560530494339251850\ldots.$$

For more details see [59]. Very recently, L. Bétermin [25] found a surprising connection between the problem of minimising a planar "Coulombian renormalised energy" derived from the Ginzburg-Landau model of superconductivity introduced by E. Sandier and S. Serfaty in [178] (also see the survey [183]) and the discrete logarithmic energy problem on $\mathbb{S}^2$. The preliminary results are: (i) the asymptotics of the minimal logarithmic energy on $\mathbb{S}^2$ has a term of order N, (ii) whose constant is bounded from above by $C_{\log,2}$ given above, and (iii) this constant equals $C_{\log,2}$ if and only if a certain triangular lattice of density 1 (called "Abrikosov" triangular lattice $\mathbb{Z} + e^{i\pi/3}\mathbb{Z}$, properly scaled, see [183]) is the minimiser of the Coulombian renormalised energy.

3.7. Numerical integration and discrepancy from the energy point of view

The reproducing kernel Hilbert space approach (see Section 2) enables us to write the squared worst-case error as

$$\frac{1}{N^2} \sum_{i=1}^{N} \sum_{j=1}^{N} K(\mathbf{x}_i, \mathbf{x}_j) - \int_{\mathbb{S}^d} \int_{\mathbb{S}^d} K(\mathbf{x}, \mathbf{y}) \, \mathrm{d}\sigma_d(\mathbf{x}) \mathrm{d}\sigma_d(\mathbf{y}),$$

which can be interpreted as K-energy of the node set $\{\mathbf{x}_1, \ldots, \mathbf{x}_N\}$ of the QMC method. Here, the energy kernel is a reproducing kernel for $H^s(\mathbb{S}^d)$. Optimal node sets are solutions of the minimal energy problem for this kernel. In [48] and [49] it is shown how the distance kernel

$$K_{\mathrm{gd}}^{(s)}(\mathbf{x}, \mathbf{y}) := 2W_{d-2s}(\mathbb{S}^d) - \|\mathbf{x} - \mathbf{y}\|^{2s-d}, \qquad \mathbf{x}, \mathbf{y} \in \mathbb{S}^d, \tag{3.19}$$

arises in a natural way as a reproducing kernel for $H^s(\mathbb{S}^d)$ for s in $(d/2, d/2 + 1)$. Evidently, there is a close connection between finding optimal QMC nodes and the problem of maximising the sum of all pairwise distances taken to the power $2s - d$. Wagner's bounds (see [199, 200]) yield that a sequence of N-point maximisers of such a generalised sum of distances is a QMC design sequence for $H^s(\mathbb{S}^d)$, $d/2 < s < d/2 + 1$; see [61]. Explicit constructions for QMC design sequences are not known. Based on numerical evidence it is conjectured in [1, 47] that a $(0, 2)$-sequence (a special digital net sequence in the sense of [88]) or a Fibonacci lattice (also see [88]) in the square $[0, 1)^2$ mapped to the 2-sphere via an *area-preserving* map (*Lambert azimuthal equal-area projection*) will be a QMC design

sequence for $H^{3/2}(\mathbb{S}^2)$. For $s \geq d/2 + 1$, the space $H^s(\mathbb{S}^d)$ can also be provided with a reproducing kernel which is essentially a distance kernel with power $2s - d$. In order to ensure that the reproducing kernel is positive definite (in the sense of Schoenberg [181]) a polynomial correction term is needed when $s - d/2$ is not a positive integer.[26] [27] Such a correction term is annihilated when the search for optimal QMC designs for $H^s(\mathbb{S}^d)$, $s \in (d/2 + L, d/2 + 1 + L)$, is restricted to spherical L-designs. In that case it suffices to minimise the energy functional (see [61])

$$[\mathrm{wce}_{H^s}(X_{N,L})]^2 = \frac{1}{N^2} \sum_{i=1}^{N} \sum_{j=1}^{N} (-1)^{L+1} \|\mathbf{x}_{i,L} - \mathbf{x}_{j,L}\|^{2s-d} - (-1)^{L+1} W_{d-2s}(\mathbb{S}^d)$$

subject to the condition that the node set $X_{N,L} = \{\mathbf{x}_{1,L}, \ldots, \mathbf{x}_{N,L}\} \subset \mathbb{S}^d$ is a spherical L-design. Note that the fixed L (once s is fixed) is "small" in the following sense: a QMC method with a more regular (in the sense of spherical designs) node set is more suitable for integrating functions from a smoother function space $H^s(\mathbb{S}^d)$. But regularity beyond a critical order will not improve the worst-case error bound; this should be compared with the optimal order (2.38). We remark that the Cui and Freeden kernel [82]

$$\mathrm{K}_{\mathrm{CF}}(\mathbf{x}, \mathbf{y}) := 2 - 2 \log \left(1 + \frac{1}{2} \|\mathbf{x} - \mathbf{y}\| \right), \qquad \mathbf{x}, \mathbf{y} \in \mathbb{S}^2,$$

which was used to define a "generalised discrepancy" to measure uniform distribution of point set sequences, can be interpreted as reproducing kernel for $H^{3/2}(\mathbb{S}^2)$ as observed in [187] and the minimising K_{CF}-energy point configurations give rise to a QMC design sequence for $H^{3/2}(\mathbb{S}^2)$. Recently, C. Choirat and R. Seri [69, 70] derived the analogous kernel for d-spheres. The corresponding minimising configurations then form QMC design sequences for $H^{(d+1)/2}(\mathbb{S}^d)$. A curious observation is that when $H^s(\mathbb{S}^d)$, $d/2 < s < d/2+1$, is provided with the reproducing kernel $\mathrm{K}_{\mathrm{gd}}^{(s)}$ from (3.19), a limit process yields that (see [62])

$$\lim_{s \to (d/2)^+} \frac{[\mathrm{wce}_{H^s}(X_N)]^2 - \frac{1}{N}}{2s - d} = \frac{1}{N^2} \sum_{\substack{i=1 \\ i \neq j}}^{N} \sum_{j=1}^{N} \log \frac{1}{\|\mathbf{x}_i - \mathbf{x}_j\|} - V_{\log}(\mathbb{S}^d)$$

for any N-point set $X_N = \{\mathbf{x}_1, \ldots, \mathbf{x}_N\} \subset \mathbb{S}^d$. This suggests that the function space $H^{d/2}(\mathbb{S}^d)$ (which is not a reproducing kernel Hilbert space) is paired with the logarithmic kernel and the logarithmic energy of an N-point set can be understood as a limit of worst-case errors in the above sense. This pairing can be extended to $s < d/2$ and the open

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[26]A logarithm of the distance appears in the reproducing kernel when $s - d/2$ is a positive integer.

[27]When defining the function space as Bessel potential space, then no correction terms are needed. In the Hilbert space setting ($p = 2$), the worst-case error is given as a Bessel-energy. The Bessel kernel on the sphere, however, has a series expansion in spherical harmonics without a convenient closed form representation. For general $p > 1$ the worst-case error has an integral representation; see [51] and also [42].

question is how to define integration and error of integration so that the corresponding Riesz $(2s − d)$-energy is a meaningful measure for this error.

Due to J. Beck [18], the spherical cap discrepancy $D(X_N)$ given in (2.16) of any $X_N \subset \mathbb{S}^d$ is bounded like $\gg N^{-1/2-1/(2d)}$ and, by appealing to a probabilistic argument, there are $X_N \subset \mathbb{S}^d$ with

$$D(X_N) \ll N^{-1/2-1/(2d)} \sqrt{\log N}. \tag{3.20}$$

A sequence of point sets with this property is called a *low-discrepancy sequence on $\mathbb{S}^d$*. It is shown in [61] that a low-discrepancy sequence (X_N^{LD}) on $\mathbb{S}^d$ is *almost* a QMC design sequence for $H^{(d+1)/2}(\mathbb{S}^d)$ in the sense that for sufficiently large N,

$$c\, N^{-(d+1)/(2d)} \leq \mathrm{wce}_{H^{(d+1)/2}}(X_N^{\mathrm{LD}}) \leq C\, N^{-(d+1)/(2d)} \sqrt{\log N}. \tag{3.21}$$

One of the deep unresolved questions is if the logarithmic term in (3.20) arising from a probabilistic argument can be removed. It is also unknown how to construct a sequence of N-point sets explicitly with spherical cap discrepancy decaying like $N^{-1/2-1/(2d)} \sqrt{\log N}$. For the 2-sphere, the construction of A. Lubotzky, R. Phillips, and P. Sarnak [149, 150] satisfies the estimate $D(X_N^{\mathrm{LPS}}) \ll (\log N)^{2/3} N^{-1/3}$ with numerical evidence indicating a convergence rate of $\mathcal{O}(N^{-1/2})$, whereas the *spherical Fibonacci lattice point sets* of [1] obey the estimate $D(X_N^{ABD}) \leq 44\sqrt{8}\, N^{-1/2}$ with numerical results showing a convergence rate of $\mathcal{O}((\log N)^c\, N^{-3/4})$ for some $1/2 \leq c \leq 1$. The *typical* spherical cap discrepancy of N i.i.d. uniformly distributed random points on $\mathbb{S}^2$ is of exact order $N^{-1/2}$, see [1]. Surprisingly, minimal Coulomb energy points on $\mathbb{S}^2$ do not have low spherical cap discrepancy. J. Korevaar [135] conjectured that minimal $(d − 1)$-energy configurations on $\mathbb{S}^d$ have spherical cap discrepancy of order $N^{-1/d}$. This conjecture was proven by M. Götz [101] up to a logarithmic factor. He also gave a lower bound of order $N^{-1/2}$ for $d = 2$. It is open if Korevaar's conjecture extends to the full potential-theoretic regime. On the basis of the Poppy-seed Bagel theorem, one could conjecture that minimal s-energy points for hypersingular s should have small spherical cap discrepancy. The only result obtained so far is the very weak order $\sqrt{\log \log N}/\log N$ result in [83] for the boundary case $s = d$. The proof employs a smoothed Riesz energy functional.

Stolarsky's invariance principle (see K. B. Stolarsky [191]) states that the sum of all mutual distances (a Riesz energy with Riesz parameter $−1$) and the spherical cap L^2-discrepancy

$$D_{L^2}(X_N) := \left(\int_0^\pi \int_{\mathbb{S}^d} \left| \frac{1}{N} \sum_{\mathbf{y} \in X_N} \mathbb{1}_{\mathbf{y} \in C(\mathbf{z}, \theta)} - \sigma_d(C(\mathbf{z}, \theta)) \right|^2 \mathrm{d}\sigma_d(\mathbf{z}) \sin\theta \mathrm{d}\theta \right)^{1/2}$$

is constant regardless of the choice of the node set X_N on $\mathbb{S}^d$; *i.e.*,

$$\frac{1}{N^2} \sum_{i=1}^N \sum_{j=1}^N \|\mathbf{x}_i - \mathbf{x}_j\| + \frac{1}{C_d} \left[D_{L^2}(X_N) \right]^2 = \int_{\mathbb{S}^d} \int_{\mathbb{S}^d} \|\mathbf{x} - \mathbf{y}\| \, \mathrm{d}\sigma_d(\mathbf{x}) \, \mathrm{d}\sigma_d(\mathbf{y}).$$

This principle connects in a very direct way the three areas optimal energy (maximising the sum of distances on $\mathbb{S}^d$), uniform distribution (spherical cap L^2-discrepancy), and numerical integration with QMC methods for functions on $\mathbb{S}^d$. Rearranging terms,

$$\frac{1}{\sqrt{C_d}} D_{L^2}(X_N) = \sqrt{\int_{\mathbb{S}^d} \int_{\mathbb{S}^d} \|\mathbf{x} - \mathbf{y}\| \, d\sigma_d(\mathbf{x}) \, d\sigma_d(\mathbf{y}) - \frac{1}{N^2} \sum_{i=1}^{N} \sum_{j=1}^{N} \|\mathbf{x}_i - \mathbf{x}_j\|},$$

one obtains a convenient way of computing the discrepancy $D_{L^2}(X_N)$. It is shown in [49] that the right-hand side represents the worst-case error of a QMC method with node set $\{\mathbf{x}_1, \ldots, \mathbf{x}_N\} \subset \mathbb{S}^d$ for functions in $H^s(\mathbb{S}^d)$ provided with the distance kernel $\mathrm{K}_{\mathrm{gd}}^{(s)}$ given in (3.19) with $s = (d+1)/2$. Moreover, in this setting, the spherical cap L^2-discrepancy can be interpreted as worst-case error and *vice versa*. A conjecture for the asymptotic expansion (as $N \to \infty$) of the minimal spherical cap L^2-discrepancy that is based on the fundamental conjecture for the minimal Riesz s-energy and the principle of analytic continuation is proposed in [45]. The paper [48] gives an extension of Stolarsky's invariance principle to general powers of the distance (raised to the power $2s - d$) involving the generalised spherical L_2-discrepancy

$$\int_0^{\pi} \int_{\mathbb{S}^d} |\mathcal{D}_{X_N,\beta}(\mathbf{z}; \cos\theta)|^2 \, d\sigma_d(\mathbf{z}) \sin\theta d\theta,$$

for the local discrepancy function (with smoothness index $s = \beta + (d-1)/2$)

$$\mathcal{D}_{X_N,\beta}(\mathbf{z}; t) := \frac{1}{N} \sum_{\mathbf{x} \in X_N} (\mathbf{x} \cdot \mathbf{z} - t)_+^{\beta-1} - \int_{\mathbb{S}^d} (\mathbf{y} \cdot \mathbf{z} - t)_+^{\beta-1} \, d\sigma_d(\mathbf{y}).$$

The paper [50] considers the discrepancy with respect to truncated spherical cones that are anchored at infinity and extends Stolarsky's invariance principle to this setting. In [106] further connections between energy and discrepancy are discussed. We also mention [146] which considers asymptotically uniformly distributed points with an upper bound on the spherical cap discrepancy and a lower bound on the separation. The sphere case motivated [84, 85].

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