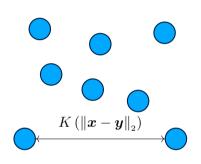
## General Kernel Spectral Methods for Equilibrium Measures An MMSC Dissertation Project



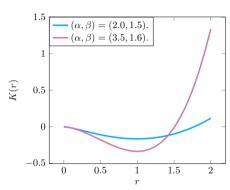
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▶ Find the Equilibrium Distribution  $\rho(x)$  of a Many-Particle-System.



N=8 particles interacting with one another through the potential K(r).



Plot of attractive-repulsive potential functions  $K(r) = \frac{r^{\alpha}}{\alpha} - \frac{r^{\beta}}{\beta}$  for different  $\alpha, \beta$ .



► Interactions through (power-law) Attraction-Repulsion Potentials¹

$$K(r) = \frac{r^{\alpha}}{\alpha} - \frac{r^{\beta}}{\beta}$$
 with parameters  $\alpha, \beta \in \mathbb{R} \setminus \{0\}$ .

▶ Each particle i = 1, ..., N at position  $x_i \in \mathbb{R}^d$  and time  $t \in \mathbb{R}^+$  follows

$$\frac{\mathrm{d}^{2}\boldsymbol{x_{i}}}{\mathrm{d}t^{2}} = f\left(\left\|\frac{\mathrm{d}\boldsymbol{x_{i}}}{\mathrm{d}t}\right\|_{2}\right) \frac{\mathrm{d}\boldsymbol{x_{i}}}{\mathrm{d}t} - \frac{1}{N} \sum_{i=1}^{N} \sum_{j\neq i}^{N} \nabla K\left(\left\|\boldsymbol{x_{i}} - \boldsymbol{x_{j}}\right\|_{2}\right),$$

for reference see, for example, [1, 2]. For now, we only consider the case without an external potential V(x).

<sup>&</sup>lt;sup>1</sup>If the repulsive term is stronger (so  $\beta > \alpha$ ), there is no equilibium distribution as particles simply continue repelling each other out to infinity.



The positional distribution approached by N=250 particles.



The total potential energy of an N-particle system is then given by

$$E = \sum_{i=1}^{N} \sum_{j=1, j \neq i}^{N} K(\|\boldsymbol{x}_{i} - \boldsymbol{x}_{j}\|_{2}),$$

which, in the continuous limit as  $N \to \infty$ , becomes

$$E = rac{1}{2} \iint K\left( \left\| oldsymbol{x} - oldsymbol{y} 
ight\|_2 
ight) \, \mathrm{d} 
ho(oldsymbol{x}) \, \mathrm{d} 
ho(oldsymbol{y}) \, ,$$

where  $d\rho = \rho(x)dx$  is a measure (the equilibrium distribution) chosen such that

$$M = \int \mathrm{d}\rho = \int_{\mathrm{supp}(a)} \rho(\mathbf{x}) \, \mathrm{d}\mathbf{x} = 1.$$



▶ In order to find  $\rho(x)$ , we consider the following ansatz

$$\rho(\boldsymbol{x}) = \sum_{n=0}^{\infty} c_n P_n^{(a,b)}(\boldsymbol{x}), \quad c_n \in \mathbb{R},$$

with which we construct a spectral method for the numerical solution of the above integral equation.

- ▶ Minimization routine of E over coefficients in  $\rho$ , as a subroutine of outer minimisation over the bounds of the box (simpler case: use [-r, r],  $r \in \mathbb{R}^+$ ).
- ▶ By construction, we find that we do not need an iterative approach for the inner optimisation routine.
- ► The outer minimisation can be performed using known methods from continuous optimisation.



Jacobi polynomials  $P_n^{(a,b)}(x)$  are orthogonal on [-1,1] w.r.t. the weight function

$$w^{(a,b)}(x) = (1-x)^a (1+x)^b,$$

so they satisfy

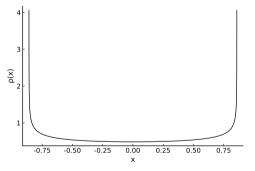
$$\int_{-1}^{1} (1-x)^{a} (1+x)^{b} P_{n}^{(a,b)} P_{m}^{(a,b)} dx = \frac{2^{a+b+1} \Gamma(a+n+1) \Gamma(b+n+1)}{n! (a+b+2n+1) \Gamma(a+b+n+1)} \delta_{n,m},$$

with a, b > -1, which uniquely determines  $P_n^{(a,b)}(x)$ . The special case of a = b corresponds to the ultraspherical or Gegenbauer polynomials, while the case a = b = 0 corresponds to the Legendre polynomials [3].

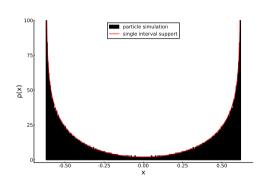
▶ This basis yields a **sparse**, and in particular, **banded** operator.



Using the spectral method, we try to immediately solve for the distribution function in the continuous limit (here, d = 1 dimension).



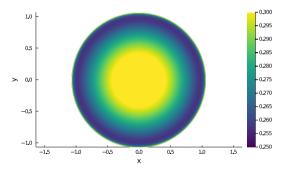
Equilibrium distribution  $(\alpha, \beta) = (2, 1.5)$  [1].



Equilib. distribution  $(\alpha, \beta) = (3.5, 1.6)$  [1].



The same is possible for d=2 or more dimensions.



The radially symmetric equilibrium distribution  $\rho(x)$  for  $\alpha = 1.2$ ,  $\beta = 0.1993$  in d = 2 dimensions, obtained using a spectral method [2].



## In this MMSC dissertation project, we will:

- ▶ Establish theory behind spectral methods in the Jacobi basis.
- $\blacktriangleright$  Use kernel expansions to construct an equilibrium measure method on [-1,1].
- ightharpoonup Numerically solve for  $\rho(x)$  using a fully implemented spectral solver<sup>2</sup>.
- ightharpoonup Implement a particle simulator for the same kernel K (mostly done).
- ▶ Compare results of the two methods and analytic solutions of special cases [4].

## Further goals include:

- ightharpoonup Considering the d-dimensional unit ball domain (d > 1).
- ightharpoonup Optionally: Considering external potentials V(x).
- ▶ Optionally: Extending to l-Morse potentials  $K(r) = C_1 e^{r/l_1} C e^{r/l}$ .



Questions?



- [1] Timon S. Gutleb, José A. Carrillo and Sheehan Olver. 'Computing equilibrium measures with power law kernels'. In: *Math. Comput.* 91.337 (Sept. 2022), pp. 2247–2281. ISSN: 0025-5718. DOI: 10.1090/mcom/3740.
- [2] Timon S. Gutleb, José A. Carrillo and Sheehan Olver. 'Computation of Power Law Equilibrium Measures on Balls of Arbitrary Dimension'. In: Constr. Approx. (Dec. 2022), pp. 1–46. ISSN: 1432-0940. DOI: 10.1007/s00365-022-09606-0.
- [3] F.W.J. Olver, A.B.O. Daalhuis, D.W. Lozier, B.I. Schneider, R.F. Boisvert, C.W. Clark, B.R. Miller and B. V. Saunders (eds.) NIST Digital Library of Mathematical Functions. http://dlmf.nist.gov. Dec. 2018. (Visited on 11/11/2020).
- [4] J. A. Carillo, Y. Huang and S. Martin. 'Explicit flock solutions for Quasi-Morse potentials'. In: European Journal of Applied Mathematics 25.5 (2014), pp. 553–578. DOI: 10.1017/S0956792514000126.

