

Start

Generate Density  
Functional  
Theory (DFT)  
DataFit  
Potential/Force

MD simulation

Calculate Forces

Calculate Next  
MD State $t_{\text{end}}?$ 

End

# MDP.jl the **julia** library of Molecular Dynamics (MD) Potentials

- ❖ Provide a comprehensive library of coupled Empirical and Machine Learning (ML) potentials

Fit DFT forces to  $\mathbf{F} = \mathbf{F}_{\text{Empirical}} + \mathbf{F}_{\text{Machine Learning}}$

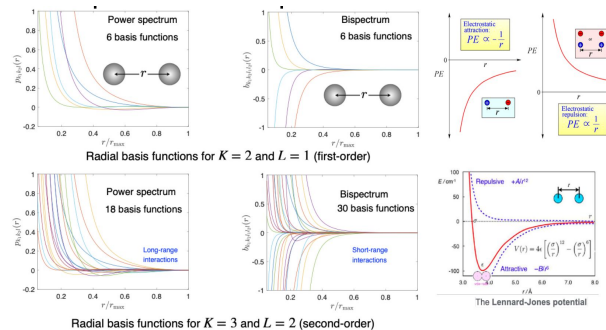
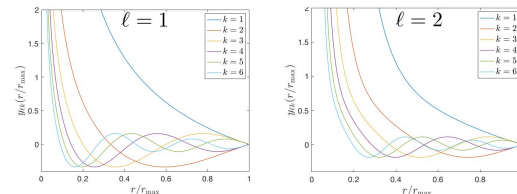
where  $\mathbf{F}_{\text{Machine Learning}} = \mathbf{F}_{\text{Power Spectrum}} + \mathbf{F}_{\text{Bispectrum}}$

- ❖ Implement automatic code generator for a wide variety of empirical potentials
- ❖ Quantify uncertainties for the trained potentials
- ❖ Increase MD simulation performance by implementing and auto-tuning new algorithms for force calculation
- ❖ Target both CPUs and GPUs
- ❖ Open Source

*“Providing fast and accurate potentials for classical MD simulations on exascale supercomputers.”*

The descriptors are based on products of spherical harmonics  $Y_{lm}(\theta, \phi)$  and spherical Bessel functions  $y_{lk}(r)$

$$u_{klm}(r) = y_{lk}(r) Y_{lm}(\theta, \phi), \quad 1 \leq k \leq K, 0 \leq l \leq L, -l \leq m \leq l$$



CESMIX

