

Computational Quantum Field Theory (CQT)

Prof. Dr. Wolfhard Janke

Institut für Theoretische Physik

wolfhard.janke@itp.uni-leipzig.de
<http://www.physik.uni-leipzig.de/cqt.html>

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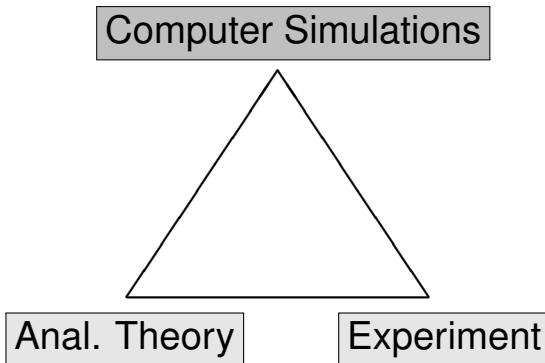
Methodologies

Anal. Theory

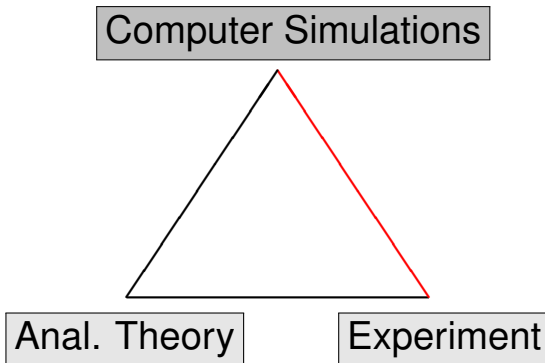
Experiment



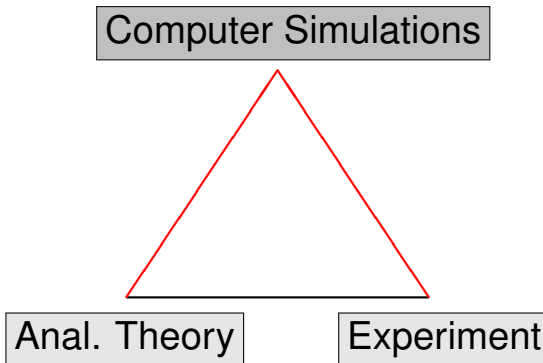
Methodologies



Methodologies



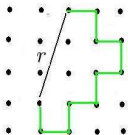
Methodologies



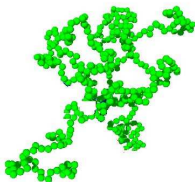
Advanced **Monte Carlo (MC)** and Molecular Dynamics (MD)
computer simulations



Polymers as Self-Avoiding Random Walks (SAWs)



N steps



$N + 1$ monomers

- SAW = random walk (RW) that is *not* allowed to cross itself
- Average end-to-end distance $\langle R_{ee} \rangle \propto N^{\nu_{\text{SAW}}}$ with “critical” exponent $\nu_{\text{SAW}} \approx \frac{3}{d+2} \geq 1/2 = \nu_{\text{RW}}$,
- Number of polymer (SAW) conformation $Z_N = c \mu^N N^{\gamma_{\text{SAW}} - 1}$

Polymer Models in the Continuum

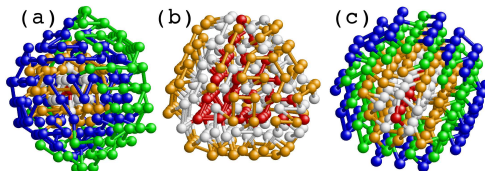
Lennard-Jones (LJ) + bond potential:

$$E = \frac{1}{2} \sum_{\substack{i,j=1 \\ i \neq j}}^N 4\epsilon [(\sigma/r_{ij})^{12} - (\sigma/r_{ij})^6] + \sum_{i=1}^{N-1} V_{\text{bond}}(r_{ii+1})$$

with

$$V_{\text{bond}}(r_{ii+1}) = -\frac{K}{2} R^2 \ln\{1 - [(r_{ii+1} - r_0)/R]^2\}.$$

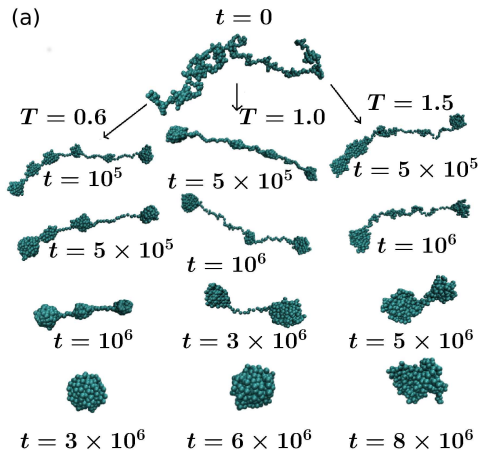
Ground-state “crystals”, similar to atomic LJ clusters:



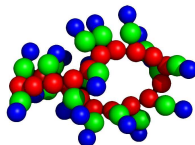
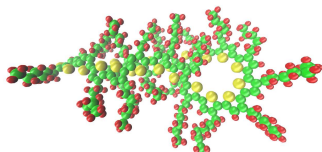
$N = 234$: icosahedral tetrahedral decahedral



Polymer Collapse Kinetics



Polymers from All-Atom to Coarse-Grained Models



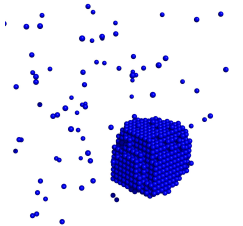
P3HT: Poly(3-hexylthiophene)

Adsorption of P3HT on gold surface in ultrahigh vacuum:
Computer simulations in comparison with experiments done at
University Halle

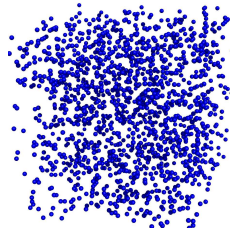


Polymer Aggregation versus Particle Condensation

Condensed



Evaporated



Balancing interface (free) energy vs fluctuations
energy vs entropy



Further Research Topics

- Disordered systems (spin glasses, long-range correlated disorder, ...)
- Method development (population annealing, long-range interactions, ...)
- Active matter systems
- Machine learning



Perspectives

- DFG Sonderforschungsbereich/Transregio SFB/TRR 102 *Polymers under Multiple Constraints: Restricted and Controlled Molecular Order and Mobility* with integrated Research Training Group (iRTG) (with Univ. Halle)
- German-French PhD College (exchange programme with Nancy, Coventry and Lviv)
- Graduate School *BuildMoNa – Building with Molecules and Nanoobjects* (with Chemistry, Biochemistry)
- International Max Planck Research School (IMPRS) *Mathematics in the Sciences* (with Math., MPI MIS)
- Cooperation with PhD school in Krakow
- EU COST action EUTOPIA



Infos

- **CQT Homepage:**
<http://www.physik.uni-leipzig.de/cqt.html>
- **CQT Report 2017:**
<http://www.physik.uni-leipzig.de/~janke/research/reports.html>
- **19th International Workshop on *New Developments in Computational Physics* “CompPhys18”:**
29 November – 01 December 2018
<http://www.physik.uni-leipzig.de/~janke/CompPhys18>
→ Programme/Abstracts of Talks

