# Predictive Compression Dynamics: A Methodological Framework for Computable Information-Motivated Modeling

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#### Abstract

We present  $Predictive\ Compression\ Dynamics\ (PCD)$ , a methodological recipe for constructing computable, local functionals  $\Phi_b$  and driving dynamics by gradient flow  $\dot{x} = -\nabla \Phi_b(x)$  with preregistered parameters. Two concrete instances are given: (i) a fixed-graph pair functional and (ii) a smooth compact-support kernel; both yield stable, attractive gradient terms (after calibration) and admit Lyapunov descent. These models serve as methodological demonstrations of computable, information-motivated optimization. We make the MDL split  $L_{\rm tot} = L(M) + L(D \mid M)$  explicit, give a minimal coding scheme linking  $\Phi_b$  descent to achievable  $\Delta L_{\rm tot}$ , address well-posedness (smooth-kernel variant), recommend robust integrators (BAOAB for Langevin), and provide a preregistration/model-card template and falsifiers for a chosen model instance. The goal is a reproducible toolbox for compression-driven dynamics across domains.

## 1 Positioning and Commitments

A disciplined workflow to construct and test *computable* local functionals  $\Phi_b$  whose gradients define algorithmic descent rules, with explicit preregistration (domain, discretization, kernels, parameters), numerical sanity checks, and internal falsifiers.

# 2 Operational Domain and Notation

We consider N point particles with positions  $x_i \in \mathbb{R}^3$  and positive weights  $m_i$ . Computations use finite precision: lattice spacing  $a_{\mathrm{grid}}$  and b bits/axis, stated a priori. The subscript b in  $\Phi_b$  denotes dependence on numerical precision (bits of representation). A global calibration constant  $G_{\mathrm{eff}} > 0$  maps dimensionless gradients to physical units if desired. Using identical  $m_i$  in both interaction and inertial terms is a modeling simplification that enforces equal accelerations by design. Vectors are written in plain type for brevity.

# 3 Model–Data Decomposition and Coding Link

Following MDL, we split description length as

$$L_{\text{tot}} = L(M) + L(D \mid M), \tag{3.1}$$

<sup>\*</sup>Jeeves is a pseudonym for an AI assistant contributing to analysis, code design, and manuscript preparation.

where L(M) encodes modeled regularities and  $L(D \mid M)$  encodes residuals given M. A decrease  $\Delta L_{\rm tot} < 0$  corresponds to realized compression. PCD treats a computable, local  $\Phi_b$  as a proxy for the achievable total codelength  $L_{\rm tot}$  itself; hence  $\dot{x} = -\nabla \Phi_b$  implements a descent in surrogate description length under the chosen model family.

Minimal explicit coding scheme. Let (i, j) range over a symmetric set of "near" pairs. A two-part code describes (i) a shared pairwise template per distance bin and (ii) residual offsets:

- Partition distances into bins  $\{B_k\}$  with centers  $r_k$ ; encode the histogram counts using an arithmetic code with probability  $p_k$  proportional to frequency.
- For each pair (i, j) with  $r_{ij} \in B_k$ , encode a residual offset  $\delta r_{ij}$  relative to  $r_k$  using bounded precision.

The expected codelength per pair is

$$\ell(r_{ij}) = -\log p_k + H_{\rm res}(\delta r \mid B_k),$$

and the overall code is prefix-free, satisfying Kraft's inequality. Then

$$L_{\text{tot}} \approx \text{const} + \sum_{(i,j)} \ell(r_{ij}).$$
 (3.2)

Choosing

$$\Phi_b \propto \sum_{(i,j)} \ell(r_{ij})$$

makes  $-\nabla\Phi_b$  a gradient descent in achievable total codelength. A convenient smooth surrogate is

$$\ell(r) \approx (r^2 + a^2)^{-1/2}$$

which provides bounded curvature at r = 0 and 1/r asymptotics. Other forms may be substituted without altering the workflow.

**Proposition 3.1** (Surrogate MDL Descent). Suppose  $L_{tot} = \text{const} + \sum_{(i,j)} \ell(r_{ij})$  with  $\ell'(r) \leq 0$  and  $\Phi_b = \kappa \sum_{(i,j)} \ell(r_{ij})$  for some  $\kappa > 0$ . Then along  $\dot{x} = -\nabla \Phi_b(x)$  we have

$$\frac{d}{dt}\Phi_b(x(t)) = -\|\nabla\Phi_b(x(t))\|^2 \le 0,$$
(3.3)

with equality iff  $\nabla \Phi_b(x(t)) = 0$ .

Interpretation.  $\Phi_b$  is a computable surrogate for total codelength; its monotone decrease under  $\dot{x} = -\nabla \Phi_b$  represents achievable compression within the chosen model family.

## 4 Information-Motivated Surrogates for Gradient Descent

#### 4.1 Fixed-graph functional

Let  $E \subset \{(i,j): 1 \leq i < j \leq N\}$  be a symmetric, degree-bounded edge set. Define

$$\Phi_E(x) = \sum_{(i,j)\in E} \frac{m_i m_j}{\sqrt{\|x_i - x_j\|^2 + a^2}}, \qquad a > 0.$$
(4.1)

The gradient term for element i is

$$G_i^{(E)}(x) = -\nabla_{x_i} \Phi_E(x) = -\sum_{\substack{j:\\(i,j)\in E}} m_i m_j \frac{(x_i - x_j)}{(\|x_i - x_j\|^2 + a^2)^{3/2}}.$$
 (4.2)

In the two-particle case with  $(i, j) \in E$  and  $a \to 0$ ,

$$G_i^{(E)} \to -m_i m_j \frac{x_i - x_j}{\|x_i - x_j\|^3},$$
 (4.3)

i.e. an attractive inverse-square form along the inter-particle direction.

#### 4.2 Smooth-kernel functional

To avoid neighbor-set discontinuities, choose a compactly supported,  $C^1$  radial kernel  $K_{\sigma}$ :  $[0,\infty) \to \mathbb{R}_{\geq 0}$  with support  $\subset [0,R\sigma]$ . Define

$$\Phi_K(x) = \sum_{i < j} m_i m_j K_{\sigma}(\|x_i - x_j\|), \tag{4.4}$$

so  $G_i^{(K)}(x) = -\nabla_{x_i}\Phi_K(x)$  is continuous and locally Lipschitz off collisions. If  $K_{\sigma}(r) \sim (r^2 + a^2)^{-1/2}$  near r = 0, one recovers the regularized two-particle form (4.3).

## 5 Algorithmic Dynamics and Integrators

We preregister all parameters  $(a_{grid}, b, a, \sigma, \Delta t, m_i, \gamma, T, seeds)$ .

**Lemma 5.1** (Compression-Rate Identity). Under  $\dot{x} = -\nabla \Phi_b(x)$  the instantaneous surrogate codelength rate is

$$\dot{\Phi}_b(t) = -\|\nabla \Phi_b(x(t))\|^2 \le 0. \tag{5.1}$$

Hence  $\Phi_b$  is a Lyapunov function and its monotone decrease represents achievable compression under the model.

Deterministic gradient flow. Explicit Euler:

$$x_i^{(t+\Delta t)} = x_i^{(t)} + \Delta t G_i(x^{(t)}), \quad G_i \in \{G_i^{(E)}, G_i^{(K)}\}.$$
(5.2)

For stability, use adaptive  $\Delta t$  or semi-implicit variants.

Stochastic descent (BAOAB recommended).

$$m_i \ddot{x}_i = G_i(x) - \gamma \dot{x}_i + \xi_i(t), \quad \langle \xi_i(t) \xi_i(t') \rangle = 2\gamma k_B T \, \delta_{ij} \delta(t - t').$$
 (5.3)

We recommend the BAOAB integrator with reported weak/strong orders.

#### 6 Sanity Checks

With  $a \to 0$  and a single pair, (4.3) holds (after calibration  $G_{\text{eff}}$ ). For  $r \gg a$ ,

$$\frac{r}{(r^2+a^2)^{3/2}} = \frac{1}{r^2} \left( 1 - \frac{3a^2}{2r^2} + O\left(\frac{a^4}{r^4}\right) \right),\tag{6.1}$$

SO

$$||G_i^{(E)}|| = m_i m_j \frac{r}{(r^2 + a^2)^{3/2}} \approx m_i m_j \frac{1}{r^2} \left(1 - \frac{3a^2}{2r^2}\right).$$
 (6.2)

These expansions serve purely as numerical consistency checks.

# 7 Well-posedness

For a > 0 and bounded degree,  $\Phi_E \in C^1(\mathbb{R}^{3N} \setminus \{x_i = x_j\})$  and  $G^{(E)}$  is locally Lipschitz off collisions. For  $C^1$  kernels with bounded  $K'_{\sigma}$ ,  $G^{(K)}$  is continuous and locally Lipschitz. Existence and uniqueness follow by Picard–Lindelöf on compact intervals. For dynamic kNN, gradients are piecewise smooth; employ hysteresis or prefer the smooth kernel.

## 8 Preregistered Model Card (example)

**Domain.**  $a_{grid} = 10 \, \mu \text{m}, b = 16.$ 

Functional.  $\Phi_K$  with Wendland  $C^2$  kernel ( $\sigma = 0.5 \,\mathrm{mm}$ ); softening  $a = 50 \,\mathrm{\mu m}$ .

**Dynamics.** BAOAB stochastic descent with  $(m_i \equiv 1, \gamma = 0.1, T = 300 \,\mathrm{K}), \Delta t = 1 \times 10^{-3} \,\mathrm{s}.$ 

**Calibration.** Fit  $G_{\text{eff}}$  by least-squares on the slope of  $||G_i^{(E)}||$  versus  $r^{-2}$  across sampled separations in a dilute two-point sandbox at  $r \gg a$ ; hold fixed thereafter.

Sanity checks. Verify (4.3) and the far-field expansion; report seeds and residuals.

#### 9 Falsifiers for a Chosen Instance

Given fixed  $(\Phi_b, \text{params})$ , declare the instance falsified if:

- (F1) Two-point trajectories disagree with the calibrated reference form beyond numerical error.
- (F2) Smooth-kernel vs fixed-graph variants differ systematically at small r beyond topology effects.
- (F3) The surrogate  $\Phi_b$  correlates poorly with *out-of-sample* compression of generated data (e.g. compare  $\Phi_b$  to actual Lempel–Ziv compression of held-out pair-distance histograms rather than the in-sample surrogate).

## 10 Discussion and Scope

PCD supplies a reproducible route from *computable* information-motivated functionals to concrete algorithmic descent schemes. That simple pairwise surrogates coincide with familiar inverse-square interactions is a feature for validation, not a claim of novelty. Future work will broaden  $\Phi_b$  (e.g. learned local codes, graph Laplacians) under the same preregistration discipline.

**Application domains.** Although demonstrated on abstract particle configurations, the same workflow applies wherever local similarity drives redundancy reduction—particle-based learning objectives, swarm control, coarse-grained fluid solvers, or clustering under computational constraints. The physical units in examples (µm-mm) serve only as scale illustrations.

# 11 Context and Relation to Existing Frameworks

PCD complements algorithmic-thermodynamic and information-geometric programs by operating directly in finite-precision configuration space, with explicitly computable surrogates and preregistered parameters. It resembles force-directed graph energies and kernel particle flows such as Stein variational gradient descent (SVGD), but contributes (i) an explicit codelength linkage via  $\Phi_b$ , (ii) a preregistered model card with declared parameters and calibration, and (iii) built-in falsifiers tied to out-of-sample compression. Unlike entropic-gravity or holographic approaches, PCD makes no physical claims beyond algorithmic optimization.

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### References

#### References

- [1] C. E. Shannon, "A mathematical theory of communication," Bell Syst. Tech. J. (1948).
- [2] J. Rissanen, "Modeling by shortest data description," Automatica (1978).
- [3] L. A. Levin, "On the notion of a random sequence," Sov. Math. Dokl. (1971).
- [4] S. Amari, Information Geometry and Its Applications, Springer (2016).
- [5] R. Jordan, D. Kinderlehrer, F. Otto, "The variational formulation of the Fokker–Planck equation," SIAM J. Math. Anal. 29 (1998).
- [6] A. Caticha, Entropic Inference and the Foundations of Physics (2012).
- [7] H. Wendland, "Piecewise polynomial, positive definite and compactly supported radial functions," Adv. Comput. Math. 4 (1995) 389–396.
- [8] B. Leimkuhler, M. Matthews, "Rational construction of stochastic numerical methods for molecular sampling," *Appl. Math. Res. eXpress* (2013).
- [9] B. Leimkuhler, C. Matthews, *Molecular Dynamics*, Springer (2016).
- [10] R. C. Prim, "Shortest connection networks and some generalizations," *Bell Syst. Tech. J.* **36**, 1389–1401 (1957).
- [11] L. Hernquist, "An analytical model for spherical galaxies and bulges," ApJ **356**, 359–364 (1990).
- [12] H. C. Plummer, "On the problem of distribution in globular star clusters," MNRAS 71, 460–470 (1911).