

# Planck Area from Half-Density Normalization

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## Abstract

Half-densities are the natural “coordinate-free integrands” for composing kernels without choosing a background measure. But choosing a *universal* convention for turning half-density objects into dimensionless numerical amplitudes introduces a  $\text{length}^{d/2}$  scale. In  $d = 4$ , this is an *area*. This note sharpens the hypothesis ladder needed for the claim “half-density normalization selects a universal area scale”, and isolates a simple dimension-matching condition under which the Planck area appears without fractional powers of couplings. A gravitational anchor based on a minimal-areal-speed principle is recorded as a separate heuristic thread [RiveroAreal] [RiveroSimple].

Minimal claim (proved under stated hypotheses): half-density composition is intrinsic, but any scalar “amplitude function” representation requires extra scalarization structure, and if that scalar representative is demanded dimensionless then a  $\text{length}^{d/2}$  constant is unavoidable. Conjectural claim (additional universality hypotheses): the scalarization constant is fixed by universal couplings and becomes a universal area scale in  $D = 4$ .

## 1. Purpose and Scope

This note isolates one technical point that is often implicit in treatments of path integrals and quantum composition: the role of half-densities (and their scaling) in making composition laws coordinate-invariant *and* dimensionally well-defined.

Claims below are labeled as **Proposition** (math-precise under hypotheses) or **Heuristic** (programmatic bridge).

**Notation (dimensions).** The half-density weight  $\text{length}^{d/2}$  uses the dimension of the manifold being integrated over in the composition law (the “composition variable” dimension). In nonrelativistic time-slicing this is typically spatial dimension; in covariant/field-theory settings it is spacetime dimension. In the gravity-only sieve (Derivation PA-D1.3), read  $d$  as the spacetime dimension  $D$  if that avoids confusion: the “ $D = 4 \Rightarrow \text{area}$ ” conclusion is about spacetime  $D = 4$ , not a claim that *spatial*  $d = 4$  is privileged.

## 2. Half-Densities and Composition Kernels

Let  $M$  be a  $d$ -dimensional manifold. A (positive) density is a section of  $|\Lambda^d T^* M|$ , and a half-density is a section of  $|\Lambda^d T^* M|^{1/2}$ .

The key operational point is: when a kernel is a half-density in its integration variable, composition of kernels does not depend on an arbitrary choice of coordinate measure.

**Heuristic PA-H1.1 (Why half-densities).** If  $K_1(x, y)$  and  $K_2(y, z)$  are chosen so that their product in the intermediate variable  $y$  is a density, then  $\int_M K_1(x, y) K_2(y, z)$  is coordinate-invariant without fixing a preferred  $dy$ . This is the structural reason why kernel composition in the path-integral formalism is coordinate-invariant.

**Derivation PA-D1.1 (Coordinate invariance of half-density pairing and composition).** In a local chart  $y = (y^1, \dots, y^d)$ , write a half-density as  $\psi(y) = \varphi(y) |dy|^{1/2}$ . Under a change of variables  $y = y(y')$ , one has  $|dy|^{1/2} = |\det(\partial y / \partial y')|^{1/2} |dy'|^{1/2}$ , so the coefficient transforms as  $\varphi'(y') = \varphi(y(y')) |\det(\partial y / \partial y')|^{1/2}$ .

Hence the product of two half-densities is a density:  $\psi_1 \psi_2 = (\varphi_1 \varphi_2) |dy|$ , and its integral is chart-independent:  $\int_M \psi_1 \psi_2$  is well-defined without choosing a background measure beyond the density bundle itself.

Kernel composition is the same mechanism: if  $K_1(x, y)$  and  $K_2(y, z)$  are half-densities in  $y$ , then  $K_1 K_2$  is a density in  $y$  and  $\int_M K_1 K_2$  is coordinate invariant.

## 3. Dimensional Analysis: Normalizing a Half-Density Requires a Scale

A density on  $M$  carries the units of  $\text{length}^d$  once physical units are assigned to coordinates. A half-density therefore carries units  $\text{length}^{d/2}$ .

**Proposition PA-P1.1 (No canonical "half-density = function" identification).** There is no canonical identification of a half-density  $\psi \in |\Lambda^d T^* M|^{1/2}$  with an ordinary scalar function  $f$  on  $M$ . Choosing such an identification is equivalent to choosing a nowhere-vanishing reference half-density  $\sigma_*$  (equivalently a positive density  $\rho_* = \sigma_*^2$ ) and writing  $\psi = f \sigma_*$ .

**Derivation PA-D1.2 (Dilation makes the  $(\text{length})^{d/2}$  weight explicit).** On  $\mathbb{R}^d$ , consider a dilation  $y \mapsto y' = ay$  with  $a > 0$ . Then  $|dy'| = a^d |dy|$ , so  $|dy'|^{1/2} = a^{d/2} |dy|^{1/2}$ . Thus even in flat space, half-densities carry an inherent  $\text{length}^{d/2}$  scaling weight.

**Derivation PA-D1.2a (Near-diagonal scaling forces the square-root Jacobian  $(\text{length})^{-d/2}$ ).** On  $M = \mathbb{R}^d$ , introduce near-diagonal coordinates  $y = x + \varepsilon v$  with  $\varepsilon > 0$ . Then  $dy = \varepsilon^d dv$ , hence  $|dy|^{1/2} = \varepsilon^{d/2} |dv|^{1/2}$ .

For a bi-half-density kernel written locally as

$$K_\varepsilon(x, y) = k_\varepsilon(x, y) |dx|^{1/2} |dy|^{1/2},$$

its pullback to  $(x, v)$  variables becomes

$$K_\varepsilon(x, x + \varepsilon v) = (\varepsilon^{d/2} k_\varepsilon(x, x + \varepsilon v)) |dx|^{1/2} |dv|^{1/2}.$$

Thus, any attempt to define a nontrivial “ $\varepsilon \rightarrow 0$ ” near-diagonal limit of kernels (the scaling step that tangent-groupoid quantization packages) inevitably produces an  $\varepsilon^{d/2}$  factor from the half-density Jacobian, and the corresponding scalar representative must be renormalized by  $\varepsilon^{-d/2}$  to stay finite. This is the same exponent as in the finite-dimensional “square-root delta” normalization: the half-density is the square root of the density Jacobian. For a canonical distributional witness carrying the same exponent, see Derivation PA-D1.2b.

**Derivation PA-D1.2b (Identity kernel: the delta bi-half-density carries the same exponent).** The identity operator on half-densities has Schwartz kernel

$$K_{\text{Id}}(x, y) = \delta^{(d)}(x - y) |dx|^{1/2} |dy|^{1/2}.$$

Under the same near-diagonal change of variables  $y = x + \varepsilon v$ , one has  $\delta^{(d)}(x - y) = \varepsilon^{-d} \delta^{(d)}(v)$  and  $|dy|^{1/2} = \varepsilon^{d/2} |dv|^{1/2}$ , hence

$$K_{\text{Id}}(x, x + \varepsilon v) = \varepsilon^{-d/2} \delta^{(d)}(v) |dx|^{1/2} |dv|^{1/2}.$$

So the universal  $\varepsilon^{-d/2}$  exponent is already encoded in the half-density identity kernel. If one scalarizes by choosing a dimensionless reference half-density  $\sigma_* := L_*^{-d/2} |dx|^{1/2}$  (for a constant length scale  $L_*$ ), then the scalar representative of the identity kernel is the **dimensionless** distribution  $k_{\text{Id}}(x, y) = L_*^d \delta^{(d)}(x - y)$ .

**Proposition PA-P1.2 (Universal \*dimensionless\* amplitudes force a  $\text{length}^{d/2}$  constant).** If one imposes the extra requirement that the scalar representative  $f$  in  $\psi = f \sigma_*$  be dimensionless in physical units, then the reference half-density  $\sigma_*$  must carry all of the  $\text{length}^{d/2}$  dimension. In particular, a *constant* (field-independent) choice of  $\sigma_*$  is equivalent to choosing a universal  $\text{length}^{d/2}$  scale.

In  $d = 4$ , this universal  $\text{length}^{d/2}$  scale is a universal *area* scale.

**Heuristic PA-H1.2 (Reciprocity claim).** Half-densities alone do not force a particular scale: the forced fact is that converting half-density objects into scalar numerical amplitudes requires extra structure (a reference half-density). The “universal area scale” claim begins only after adding two further hypotheses: 1. the reference  $\sigma_*$  is taken to be *constant* (no dependence on background metric/fields), and 2. the constant is required to be fixed by universal constants/couplings of the theory.

Under these hypotheses,  $d = 4$  is the unique dimension in which the needed  $\text{length}^{d/2}$  constant can be supplied by the gravitational coupling without fractional powers (Derivation PA-D1.3).

Derivation PA-D1.3 (Gravity-only sieve: why  $d=4$  is singled out if only  $G_d$  is used). In  $d$  spacetime dimensions, the Einstein–Hilbert action  $\frac{1}{16\pi G_d} \int d^d x \sqrt{|g|} R$  shows that (in  $c = \hbar = 1$  units) Newton’s constant has dimension  $[G_d] = \text{length}^{d-2}$ . Assume the only available dimensionful coupling used to build the universal normalization constant is  $G_d$  itself (no cosmological constant, no additional dimensionful scales), and impose PA-H2.5 in the literal “no fractional powers of  $G_d$ ” sense. Then the normalization constant has dimension  $\text{length}^{k(d-2)}$  for some integer  $k$ . Matching  $\text{length}^{d/2}$  forces  $\text{length}^{d/2} = \text{length}^{k(d-2)}$ , which holds if and only if  $d = 4$ . In that case  $G_4$  itself has dimension of area, and the corresponding area scale is the Planck area  $L_P^2 \sim \hbar G_4 / c^3$ .

Remark PA-D1.3a (Three obstruction mechanisms: why  $d=4$  is uniquely un-obstructed). The failure of the gravity-only sieve away from  $d = 4$  has three qualitatively distinct causes. For odd  $d$ , the target exponent  $d/2$  is a half-integer, while any monomial in couplings with integer length dimensions has integer total dimension — a *parity* obstruction that is categorical regardless of which integer-dimensional couplings are admitted (including  $\Lambda_d$ ). For even  $d \geq 6$ , the target  $d/2$  is an integer but  $k = d/(2(d-2)) < 1$ , so no positive-integer power of  $G_d$  alone suffices — a *magnitude* obstruction. For  $d = 2$ ,  $G_2$  is dimensionless and gravity provides no scale. Thus  $d = 4$  is not merely the solution of a single Diophantine equation; it is the unique integer dimension that evades all three obstructions simultaneously.

### 3.1 Hypotheses as Separate Knobs (What Is Forced vs Chosen)

The discussion above mixes three different kinds of statements: 1. **Geometric facts** (what half-densities are, how they compose, how they scale), 2. **Representational choices** (how one turns half-density objects into scalar numbers), 3. **Universality/selection principles** (what choices are allowed if we demand “background-free” and “built from couplings”).

To study these separately, it is useful to keep the hypotheses explicit.

**Hypothesis PA-H2.1 (Half-density formulation).** Quantum kernels are treated as bi-half-densities so that composition in intermediate variables is coordinate invariant (Section 2 and Derivation PA-D1.4).

**Hypothesis PA-H2.2 (Scalarization by a reference half-density).** To interpret half-density amplitudes as scalar numerical functions, we pick a nowhere-vanishing reference half-density  $\sigma_*$  and write  $\psi = f \sigma_*$  (Proposition PA-P1.1).

**Hypothesis PA-H2.3 (Dimensionless scalar representative).** The scalar representative  $f$  is required to be dimensionless in physical units (Proposition PA-P1.2). This forces  $\sigma_*$  to carry the full  $\text{length}^{d/2}$  weight.

**Hypothesis PA-H2.4 (Background-free constancy).** The reference  $\sigma_*$  is taken to be constant/field-independent, rather than determined by background geometry (e.g. a Riemannian volume  $|g|^{1/4}|dx|^{1/2}$ ) or by dynamical fields (e.g. a dilaton-like factor). This is the first point where a *universal constant* enters.

**Hypothesis PA-H2.5 (Analyticity / no fractional powers of couplings).**

If the universal constant is required to be built from the theory's couplings without fractional powers, then dimensional analysis becomes a *dimension sieve* rather than a tautology. This hypothesis has at least two distinct readings: 1. **Integrality (integer-exponent) reading:** the constant is a monomial in the available couplings with integer exponents (possibly allowing negative powers). Equivalently, dimension-matching becomes an integer (Diophantine) constraint on the exponents. 2. **Perturbative analyticity reading (stronger):** the constant admits a Taylor expansion around a specified weak-coupling limit, so only nonnegative integer powers appear.

To keep the branches explicit, we will refer to these as: - **Hypothesis PA-H2.5a (Integrality / monomial sieve).** The normalization constant is a monomial in admitted couplings with integer exponents. - **Hypothesis PA-H2.5b (Perturbative analyticity sieve).** The dependence is analytic at a specified weak-coupling point (hence forbids negative powers and other non-analytic dependence).

**Heuristic PA-H2.5b1 (Analyticity for dimensionful couplings needs a reference scale).** If a coupling  $g_i$  is dimensionful, “weak coupling” is only meaningful after choosing a reference scale  $\mu$  and forming a dimensionless parameter  $\hat{g}_i(\mu) := \mu^{-a_i} g_i$  (where  $[g_i] = \text{length}^{a_i}$ ). In that sense, perturbative analyticity is naturally analyticity in  $\hat{g}_i(\mu)$  near  $\hat{g}_i = 0$ . Demanding at the same time that the scalarization constant be a  $\mu$ -independent universal constant pushes one back to either: 1. an engineering-dimension monomial in admitted dimensionful couplings with nonnegative integer exponents (the sieve branch), or 2. a non-analytic RG-invariant transmutation scale (outside PA-H2.5b).

Separately, the RG “dimensional transmutation” mechanism (Heuristic PA-H2.13 / Derivation PA-D1.6a) supplies a scale *outside* both PA-H2.5a and PA-H2.5b: it is typically non-analytic in the coupling.

Derivation PA-D1.3 is the simplest gravity-only instance under the integrality reading: “use  $G_d$  without fractional powers” singles out  $d = 4$ .

**Heuristic PA-H2.6 (Where “special dimensions” can appear).** Special dimensions do not come from half-densities alone (Hypothesis PA-H2.1). They appear only after adding a selection principle like PA-H2.4–PA-H2.5: the requirement that the scalarization choice be universal, background-free, and coupling-built in a restricted (e.g. analytic) way.

**Remark PA-H2.6b (Minimal set and non-bookkeeping content).** The minimal hypothesis set for “half-density normalization  $\Rightarrow$  Planck area in  $d = 4$ ” is: PA-H2.1 (composition forces the  $d/2$  exponent — Derivation PA-D1.4a, not

a convention), PA-H2.4 (background-free scalarization — a physical principle, not bookkeeping: dropping it lets the metric-derived  $|g|^{1/4}|dx|^{1/2}$  work in any  $d$  without a new scale), PA-H2.5a (integrality sieve — a UV hypothesis: dropping it allows transmutation to supply scales in any  $d$ ), and the gravity-only coupling restriction. Each ingredient is necessary: removing any one defeats the conclusion. The Diophantine equation  $k(d-2) = d/2$  is the algebraic core, but its ingredients are not arbitrary — composition, background-freeness, and integrality each have independent physical or structural motivation.

**Heuristic PA-H2.6a (Independent D=4 filter: conformal scalarization-gauge simplicity).** The “dimension sieve” (PA-H2.5) is not the only way a special dimension can appear once one insists on scalar representatives. A different kind of filter comes from asking for *simplicity of how kinetic operators depend on the scalarization gauge*.

In the QFT-facing half-density calculus, a background metric supplies a reference half-density  $|g|^{1/4}|dx|^{1/2}$ , and the scalar Laplacian  $\Delta_g$  conjugates to an operator on half-densities

$$\tilde{\Delta} := |g|^{1/4} \Delta_g |g|^{-1/4}.$$

In a conformal background  $g_{\mu\nu} = e^{2\sigma(x)} \delta_{\mu\nu}$  in spacetime dimension  $D$ , a direct computation gives

$$\tilde{\Delta}\psi = e^{-2\sigma} \left( \partial^2 \psi - 2 \partial\sigma \cdot \partial\psi - \frac{D}{2} (\partial^2 \sigma) \psi + \frac{D(4-D)}{4} (\partial\sigma)^2 \psi \right).$$

Thus the quadratic-gradient term  $(\partial\sigma)^2 \psi$  cancels at  $D = 4$  (within the conformal ansatz). If one adopts the extra criterion that scalarization-gauge changes should not generate such quadratic-gradient “potentials” in the half-density kinetic operator, then  $D = 4$  is singled out by *operator simplicity* rather than by coupling-dimension matching.

This filter is independent of PA-H2.5 and, by itself, does not supply a length scale; it is recorded only as an additional “special dimension” candidate knob to compare against the scale-sieve hypotheses.

### 3.2 What Changes When a Hypothesis Is Relaxed?

This subsection records the main “branches” that need separate study.

1. **Drop PA-H2.3 (allow dimensionful  $f$ ).** Then no universal length <sup>$d/2$</sup>  constant is forced; the dimensional weight can be carried by the scalar representative itself (as in the usual statement “wavefunctions have dimension length <sup>$-d/2$</sup> ”).
2. **Drop PA-H2.4 (allow background geometry).** Then  $\sigma_*$  can be chosen from a metric (or other structure), and the “universal constant” is replaced by background-dependent normalization.
3. **Drop PA-H2.5a/PA-H2.5b (allow fractional powers or other non-analytic dependence).** Then in any  $d > 2$  one can build a length <sup>$d/2$</sup>

constant from gravity via  $G_d^{d/(2(d-2))}$  (in  $c = \hbar = 1$  units), so  $d = 4$  is no longer singled out; instead,  $d = 4$  is simply the unique case where the exponent is an integer.

4. **Change “which coupling supplies the scale”.** Using other dimensional couplings (cosmological constant, string tension, gauge couplings in various dimensions, etc.) yields different “special-dimension” sieves. This is conceptually aligned with the observation that some dimensions are singled out by other structures (division algebras, special holonomy, supersymmetry), but those filters are separate from the half-density story and should not be conflated.
5. **Replace PA-H2.5 by transmutation (RG as the scale supplier).** If one allows RG-invariant scales generated by compatibility (dimensional transmutation), then even dimensionless couplings can supply a physical length scale (Heuristic PA-H2.13 / Derivation PA-D1.6a). This branch does not “sieve dimensions” in the same way as PA-H2.5a/PA-H2.5b; it supplies a scale by non-analytic RG invariants rather than by monomials in couplings.

**Branch summary (keep the knobs separate).** There are three distinct “scale supplier” mechanisms in play: 1. **Monomial sieve** (PA-H2.5a/PA-H2.5b): build the needed  $\text{length}^{d/2}$  constant from admitted couplings using restricted (e.g. analytic) dependence. 2. **Fractional/non-analytic coupling dependence:** allow fractional powers or other non-analytic dependence directly in couplings (dimensional analysis becomes permissive; the  $d = 4$  sieve disappears). 3. **Dimensional transmutation:** generate a scale from RG compatibility (Derivation PA-D1.6a / Example PA-E5), which is typically non-analytic in naive coupling expansions.

### 3.3 Starting with PA-H2.5a: Integrality as a Dimension Sieve

The point of PA-H2.5a/PA-H2.5b is not that dimensional analysis alone selects a unique scale (it does not), but that *restricting allowed functional dependence on couplings* can turn dimensional analysis into a selection principle.

Derivation PA-D1.6 (PA-H2.5a: integer-exponent form of “no fractional powers”). Work in  $c = \hbar = 1$  units for dimension counting. Let the available couplings  $\{g_i\}$  have length dimensions  $[g_i] = \text{length}^{a_i}$ . Under the integrality reading of PA-H2.5, the universal normalization constant is a monomial  $C = \prod_i g_i^{n_i}$  with integers  $n_i$ . Its length dimension is  $[C] = \text{length}^{\sum_i n_i a_i}$ . Requiring  $[C] = \text{length}^{d/2}$  is therefore the integer-exponent (Diophantine) condition

$$\sum_i n_i a_i = \frac{d}{2}.$$

Existence (and non-uniqueness) of solutions depends on: 1. which couplings are admitted as “universal” inputs, and 2. whether one allows negative exponents (non-analytic at zero coupling) or insists on perturbative analyticity (nonnegative exponents).

**Heuristic PA-H2.7 (Why PA-H2.5 needs a "what counts as a coupling" rule).** If one allows arbitrary redefinitions of couplings (e.g. adjoining a new symbol  $\tilde{G} = G_d^{1/(d-2)}$ ), then “no fractional powers” becomes vacuous: the forbidden root has simply been renamed as an allowed coupling. PA-H2.5 is meaningful only together with a prior criterion for admissible coupling dependence (e.g. perturbative analyticity around a distinguished limit such as  $G_d \rightarrow 0$ ).

**Heuristic PA-H2.7a (Admissible couplings: exclude scheme parameters and non-analytic reparametrizations).** To keep PA-H2.5 from collapsing into a coordinate artifact, we implicitly adopt the following convention: 1. **Admitted couplings** are the independent parameters that appear as coefficients of local operators in the UV action after fixing a canonical normalization convention for fields (so that field-rescaling freedom is not used to hide roots/powers). 2. We allow only **analytic** reparametrizations near a chosen base point (“weak coupling”), so adjoining  $\tilde{g} = g^{1/2}$  is disallowed when it is non-analytic at that base point. 3. We explicitly exclude scheme/scale conventions from the coupling set: the renormalization scale  $\mu$ , regulator cutoffs  $\Lambda$ , and finite-subtraction constants; and we also exclude the scalarization gauge scale  $L_*$  itself (since constraining  $L_*$  is the point of the ladder).

Under this convention, PA-H2.5a is best viewed as a computational proxy once coupling coordinates are fixed, while PA-H2.5b (analyticity at the base point) is the more invariant “no roots” statement. The phrase “canonical normalization of fields” in (1) is itself a convention choice; the point is that fixing such a convention makes the admissibility rule a controlled knob rather than an implicit loophole.

**Example PA-E1 (Gravity-only).** With only  $G_d$  available,  $a_1 = d - 2$  and the condition becomes  $n(d - 2) = d/2$ . For integer  $d \geq 3$ , this has a solution only at  $d = 4$  with  $n = 1$ , reproducing Derivation PA-D1.3.

**Example PA-E2 (Gravity + cosmological constant).** If one also allows the cosmological constant  $\Lambda_d$  with  $[\Lambda_d] = \text{length}^{-2}$ , then the condition becomes  $n(d - 2) - 2m = d/2$  for integers  $n, m$ . A simple family of solutions exists for  $d$  divisible by 4: take  $n = 1$  and  $m = d/4 - 1$ , so

$$C \sim G_d \Lambda_d^{d/4-1},$$

has dimension  $\text{length}^{d/2}$ . Thus, even under PA-H2.5,  $d = 4$  is not automatically unique once additional dimensionful couplings are admitted; what is special



about  $d = 4$  in this family is that it is the only case with  $m = 0$  (no need to involve  $\Lambda_d$ ).

**Example PA-E3 (Yang--Mills coupling as an alternative sieve).** In  $d$  spacetime dimensions, the Yang--Mills action is typically written as  $\frac{1}{4g_d^2} \int d^d x F_{\mu\nu} F^{\mu\nu}$ , so  $[g_d^2] = \text{length}^{d-4}$  (equivalently  $[g_d] = \text{length}^{(d-4)/2}$ ). If we (hypothetically) allow the half-density normalization constant to be a pure monomial in  $g_d$ ,  $C \sim g_d^p$  with integer  $p \geq 0$ , then

$$[C] = \text{length}^{p(d-4)/2}.$$

Matching  $[C] = \text{length}^{d/2}$  gives the integer-exponent condition

$$p(d-4) = d \implies d = \frac{4p}{p-1} = 4 + \frac{4}{p-1}.$$

Thus integer solutions occur only when  $p-1 \mid 4$ , i.e.  $p \in \{2, 3, 5\}$ , giving  $d \in \{8, 6, 5\}$  respectively.

In particular, in  $d = 4$  the gauge coupling is dimensionless and cannot by itself supply the  $\text{length}^{d/2}$  factor needed for half-density scalarization; in that case the scale must come from another dimensionful coupling (e.g. gravity) or from a non-analytic mechanism (dimensional transmutation).

**Example PA-E4 (A universal area parameter  $\backslash(\backslash\alpha\_ast\backslash)$  as a scale supplier).** Suppose the UV theory admits a universal area parameter  $\alpha_*$  with dimension  $[\alpha_*] = \text{length}^2$  (for example, in perturbative string theory  $\alpha_*$  is the familiar  $\alpha' = l_s^2$ , with string tension  $T \sim 1/\alpha'$ ). If one allows the half-density normalization constant to be built from  $\alpha_*$  alone as a monomial  $C \sim (\alpha_*)^n$  with integer  $n$ , then

$$[C] = \text{length}^{2n},$$

and matching  $[C] = \text{length}^{d/2}$  forces  $2n = d/2$ , i.e.  $d = 4n$ . So  $\alpha_*$  provides a “background-free” source of scale but does not single out  $d = 4$  on its own; it selects dimensions divisible by 4 under the strict integrality reading of PA-H2.5. In  $d = 4$  it yields directly an area scale  $C \sim \alpha_*$  (e.g.  $C \sim \alpha'$  in the string example).

**Remark PA-E4a (Emergent string tension is a transmutation-scale instance, not a UV parameter).** In confining phases one often defines an effective string tension  $\sigma$  with  $[\sigma] = \text{length}^{-2}$ , so  $\sigma^{-1}$  supplies an area scale. Operationally, for a large rectangular Wilson loop of spatial size  $R$  and Euclidean time extent  $T$ , one defines  $V(R) = -\lim_{T \rightarrow \infty} \frac{1}{T} \ln \langle W(R \times T) \rangle$ , and an area law  $\langle W(R \times T) \rangle \sim e^{-\sigma RT}$  implies  $V(R) \sim \sigma R$  and identifies  $\sigma$  [Greensite2003Confinement]. In  $d = 4$  gauge theories with dimensionless couplings, such a scale is expected to arise (when it exists) as an RG-invariant transmutation scale rather than as a new analytic monomial in the couplings.

Logically, this places “string tension as area supplier” in the PA-H2.13 branch unless one is explicitly assuming a fundamental UV area parameter  $\alpha_*$ .

**Heuristic PA-H2.12 (Link to gravity/Planck length in a UV completion).**

The gravity-only sieve (Derivation PA-D1.3) uses  $G_d$  as the unique universal coupling supplying dimension. In a UV completion where gravity is emergent from a sector with a universal area parameter  $\alpha_*$  and a dimensionless coupling  $g$ , dimensional analysis suggests

$$G_d \propto g^2 (\alpha_*)^{(d-2)/2} \times (\text{volume factors}),$$

so the Planck length/area is *derived* from  $\alpha_*$  and  $g$  rather than fundamental. In perturbative string theory, one has  $\alpha_* = \alpha'$  and  $g = g_s$  (up to compactification-volume factors). In that framing the half-density “universal area scale” could naturally be  $\alpha_*$  (or a simple function of it), while the Planck area is recovered as a consequence of how gravity emerges.

**Heuristic PA-H2.8 (What PA-H2.5 is really buying).** The value of PA-H2.5 is comparative: it distinguishes dimensions in which the needed length <sup>$d/2$</sup>  factor can be supplied by *simple* coupling dependence (integer powers of the already-present couplings), versus dimensions in which any such factor requires either (i) introducing extra scales/couplings, (ii) taking fractional powers, or (iii) invoking non-analytic mechanisms (dimensional transmutation).

**Heuristic PA-H2.13 (Dimensional transmutation as a scale-supplier).**

If one relaxes PA-H2.5’s “analytic monomial in couplings” requirement, then even a theory with only *dimensionless* couplings can generate a physical length scale through RG invariance: a running coupling  $g(\mu)$  can be traded for an RG-invariant scale  $\kappa_*$  (or  $\Lambda$ ), typically of the form  $\kappa_* \sim \mu \exp(-\text{const}/g(\mu)^2)$  or  $\kappa_* \sim \mu \exp(-\text{const}/g(\mu))$ . In that branch, the half-density scalarization scale required by PA-H2.4 can be supplied by  $\kappa_*^{-d/2}$  (or its square in  $d = 4$  as an area scale), but the scale is no longer an analytic monomial in the couplings: it is emergent and non-perturbative in the naive coupling expansion.

**Heuristic PA-H2.14 (Bookkeeping: what “ $\text{length}^{d/2}$ ” means in “ $\text{length}^{d/2}$ ”).**

The half-density weight  $\text{length}^{d/2}$  refers to the dimension of the manifold whose coordinates are integrated over in the composition law (the intermediate-variable space). In a nonrelativistic time-sliced kernel this is typically the *spatial* dimension, while for covariant/proper-time kernels one may compose over *spacetime* points. The dimension-sieve discussion using  $G_d$  treats  $d$  as the **spacetime** dimension, so any  $d = 4 \Rightarrow$  “area scale” conclusion should be read in that covariant sense unless stated otherwise.

**Derivation PA-D1.6a (RG-invariant scale from a beta function).** Let a (dimensionless) running coupling  $g(\mu)$  satisfy an RG equation  $\mu dg/d\mu = \beta(g)$  with  $\beta(g) \neq 0$  in the range of interest. Then the combination

$$\Lambda_* \equiv \mu \exp\left(-\int^{g(\mu)} \frac{dg'}{\beta(g')}\right)$$

is RG-invariant (independent of the subtraction scale  $\mu$ ), up to a finite multiplicative constant corresponding to a choice of scheme/normalization of the integral. In one-loop form  $\beta(g) = -bg^2 + O(g^3)$ , one obtains the familiar transmutation scale  $\Lambda_* \sim \mu e^{-1/(bg(\mu))} \times (\text{scheme factor})$ .

If PA-H2.3–PA-H2.4 demand a universal scalarization constant  $C$  with  $[C] = \text{length}^{d/2}$ , then any RG-invariant inverse length  $\Lambda_*$  supplies one by  $C \sim \Lambda_*^{-d/2}$ . In particular, for  $d = 4$  this produces a universal **area** scale  $C \sim \Lambda_*^{-2}$ , without requiring the scale to be an analytic monomial in couplings (so this branch sits outside PA-H2.5).

**Example PA-E5 (2D delta: transmutation yields a length scale).** In the 2D delta interaction, the contact coupling is marginal and the renormalized theory is naturally parameterized by an RG-invariant inverse length  $\kappa_*$  rather than by the bare coupling. Concretely, one finds (up to conventions) a running coupling  $g_R(\mu)$  with beta function  $\beta(g_R) \propto g_R^2$ , and the RG invariant

$$\kappa_*^2 \equiv \mu^2 \exp\left(\frac{2\pi\hbar^2}{m} \frac{1}{g_R(\mu)}\right),$$

so  $\kappa_*$  is independent of the subtraction scale  $\mu$  and sets a bound-state/scattering scale [ManuelTarrach1994PertRenQM]. This is a minimal witness that “a scale is forced by compatibility” can occur even without a dimensionful coupling, via renormalization rather than via analytic monomials.

**Remark PA-E5a (Half-density match and three-level RG hierarchy).** In the 2D delta, the transmutation scale  $\kappa_*^{-1}$  has dimension  $\text{length}^1 = \text{length}^{d/2}$  for  $d = 2$ , exactly the half-density scalarization weight required by PA-H2.3–PA-H2.4. The beta function  $\beta(g_R) \propto g_R^2$  vanishes to order 2 at the Gaussian fixed point; transmutation requires this nonlinearity — a linear beta function produces only algebraic (power-law) RG invariants without generating a new scale. Three levels of the RG hierarchy are now witnessed by explicit computations: 1. *Semigroup structure*: shared by all refinement flows, including linear beta functions (any toy ODE with a linear beta, e.g.  $\beta(a) = \frac{1}{2} - a$ , exhibits the semigroup property without transmutation). 2. *Transmutation*: requires  $\beta$  of order  $\geq 2$  at the fixed point, producing a non-analytic RG-invariant scale. Witness: this example (2D delta,  $\beta \propto g_R^2$ ). 3. *Dimension sieve* (PA-H2.5): demands the scale be an analytic monomial in couplings, selecting  $d = 4$  under the gravity-only hypothesis. Witness: Derivation PA-D1.3. Transmutation (level 2) supplies a scale in any  $d$  where a marginal coupling exists, so it does not sieve dimensions. The half-density weight  $\text{length}^{d/2}$  correctly tracks the geometric type of the resulting scale: a length in  $d = 2$ , an area in  $d = 4$ .

**Remark PA-E5b (Where in the kernel the transmutation scale acts).** The full resolvent of the 2D delta interaction factorizes via the Lippmann–Schwinger identity as  $G = G_0 + G_0 T G_0$ , where  $G_0(x, z; E)$  is the free resolvent and  $T(E)$  is the scalar  $T$ -matrix at the contact vertex. Each  $G_0$  factor carries the Van Vleck half-density weight at its endpoint (Derivation PA-D1.4), while

$T(E)$  is a scalar amplitude containing the transmutation scale  $\kappa_*$ . Thus the two ingredients play complementary roles: the Van Vleck prefactor implements the half-density *transformation law* (PA-H2.1), and the transmutation scale supplies the *scalarization constant* needed to extract dimensionless amplitudes (PA-H2.2–PA-H2.4). They are structurally independent and combine multiplicatively in the kernel; the geometric weight comes from free propagation between vertices, while the dynamical scale comes from RG invariance at the vertex.

**Remark PA-E5c (Half-density composition at the vertex mirrors the free-propagator semigroup).** In the Lippmann–Schwinger product  $G_0 T G_0$ , the half-density mechanism is the same as in PA-D1.4a: the two  $|dy|^{1/2}$  factors at the vertex pair into a density  $|dy|$ , which integrates coordinate-invariantly, while the result inherits  $|dx|^{1/2} |dz|^{1/2}$  at the endpoints. The  $T$ -matrix is a scalar at the contact vertex containing the transmutation scale  $\kappa_*$  but carrying no half-density weight. Thus the half-density structure (geometric, from free propagation) and the transmutation scale (dynamical, from RG at the vertex) enter the resolvent through independent factors and combine multiplicatively. The transmutation scale  $\kappa_*^{-1}$  has dimension  $\text{length}^{d/2}$  for  $d = 2$ , matching the half-density scalarization weight. For  $d = 4$ , the same factorization gives a scalarization constant  $\Lambda_*^{-d/2} = \Lambda_*^{-2} = \text{area}$ .

### 3.4 Running PA-H2.3: Is “Dimensionless $f$ ” Physics or Convention?

The half-density formalism (PA-H2.1) gives a canonical pairing  $\int \bar{\psi} \psi$  that does not require choosing a background measure. But when we write  $\psi = f \sigma_*$  (PA-H2.2), we are choosing a *representation* of the same object as a scalar function with respect to a chosen positive density  $\rho_* = \sigma_*^2$ .

**Proposition PA-P1.3 (Scalarization is a choice of measure, not new physics).** Choosing a reference half-density  $\sigma_*$  identifies the canonical Hilbert space of  $L^2$  half-densities on  $M$  with the scalar Hilbert space  $L^2(M, \rho_*)$ , where  $\rho_* = \sigma_*^2$ . Different choices  $\sigma_*$  yield unitarily equivalent scalar representations.

**Derivation PA-D1.7 (Change of reference half-density acts by multiplication).** Let  $\sigma_1, \sigma_2$  be nowhere-vanishing half-densities on  $M$  and set  $r := \sigma_2/\sigma_1$ , a positive scalar function. Writing the same half-density state  $\psi$  as  $\psi = f_1 \sigma_1 = f_2 \sigma_2$  gives  $f_2 = r^{-1} f_1$ . Moreover,

$$\int_M \bar{\psi} \psi = \int_M |f_1|^2 \sigma_1^2 = \int_M |f_2|^2 \sigma_2^2 = \int_M |f_2|^2 r^2 \sigma_1^2,$$

so the two scalar pictures differ by a compensating change of measure and pointwise multiplication. In particular, if  $\sigma_2 = c \sigma_1$  is a constant rescaling, then  $f_2 = c^{-1} f_1$  is the familiar global wavefunction normalization freedom.

**Heuristic PA-H2.9 (How PA-H2.3 creates a scale).** In the usual “scalar wavefunction” presentation on  $\mathbb{R}^d$ , one implicitly chooses  $\sigma_* = |dx|^{1/2}$  and

allows the scalar representative to carry dimension  $\text{length}^{-d/2}$  so that  $|\psi|^2 d^d x$  is dimensionless probability. Requiring instead that the scalar representative  $f$  be dimensionless (PA-H2.3) shifts the  $\text{length}^{-d/2}$  factor into the reference half-density:

$$\sigma_* \sim L_*^{-d/2} |dx|^{1/2},$$

so “dimensionless  $f$ ” is a convention unless the scale  $L_*^{d/2}$  is fixed by an additional universality principle (PA-H2.4–PA-H2.5).

### 3.5 Running PA-H2.4: What Does “Background-Free Constancy” Mean?

From Derivation PA-D1.7, changing the reference half-density  $\sigma_*$  by a positive function  $r(x)$  changes the scalar representative by  $f \mapsto r^{-1}f$  and changes the scalar measure by  $\rho_* \mapsto r^2 \rho_*$ . So the raw half-density formulation has a large “scalarization gauge freedom”.

**Heuristic PA-H2.10 (Constancy = no extra background function).**

If we take “background-free” in the strong sense “no additional structure beyond the manifold and the theory’s couplings”, then allowing an arbitrary non-constant  $r(x)$  would amount to introducing a new background field/function by hand. In that strong sense, the only admissible changes are constant rescalings, and choosing  $\sigma_*$  becomes a choice of a single global scale (fixed or not fixed by couplings depending on PA-H2.5).

**Derivation PA-D1.8 (Three natural families of  $\sigma_*$  and what they mean).** On a configuration space  $M$ , the common ways to choose a reference half-density are: 1. **Flat/affine choice (when available):** on  $\mathbb{R}^d$  with its affine structure, translation invariance picks  $|dx|^{1/2}$  uniquely up to a constant factor. This is “constant” in the sense of being homogeneous under translations. 2. **Metric-derived choice:** given a Riemannian/Lorentzian metric  $g$ , one can take  $\sigma_g := |g|^{1/4} |dx|^{1/2}$ , so that  $\rho_g = \sigma_g^2 = \sqrt{|g|} |dx|$  is the familiar invariant volume density. This makes the scalar representative  $f$  a genuine scalar field but makes the scalarization depend on background geometry. 3. **Field-derived (dilaton-like) choice:** given a scalar field  $\Phi$  (background or dynamical), one can take  $\sigma_\Phi := e^{-\Phi} \sigma_g$ . In the scalar picture this is a local rescaling of the measure, and it is the natural way to encode “local units” or Weyl factors.

PA-H2.4 asserts that the theory supplies (or selects) a choice of type (1) with no  $x$ -dependent factor: a fixed reference  $\sigma_*$  whose only remaining ambiguity is an overall constant scale.

**Heuristic PA-H2.11 (RG as scale dependence of scalarization).** If refinement/coarse-graining forces an  $x$ -independent but scale-dependent choice  $\sigma_*(\mu)$  (equivalently a scale-dependent constant  $L_*(\mu)$ ), then PA-H2.4 is replaced by an RG statement: the scalarization convention becomes part of the renormalization scheme (a “wavefunction renormalization” for the scalar

representative). In that case, a universal area/length scale can still appear, but typically as an RG invariant (dimensional transmutation scale) rather than as a fixed analytic monomial in couplings.

**Derivation PA-D1.8a (Running scalarization  $\sigma_*(\mu)$  is a  $Z(\mu)$  factor on scalar representatives).** Assume the intrinsic (half-density) state  $\psi$  is fixed, and consider two scalarizations related by a  $\mu$ -dependent *constant* rescaling,  $\sigma_*(\mu) = c(\mu) \sigma_0$  with  $c(\mu) > 0$ . Writing  $\psi = f(\mu) \sigma_*(\mu)$  gives  $f(\mu) = c(\mu)^{-1} f_0$ , where  $f_0$  is the scalar representative in the  $\sigma_0$  convention. Defining  $Z(\mu) := c(\mu)^2$ , this becomes the familiar multiplicative form

$$f(\mu) = Z(\mu)^{-1/2} f_0.$$

So an  $x$ -independent running of scalarization is *formally equivalent* to a wave-function renormalization factor on scalar representatives. This is bookkeeping:  $\sigma_*(\mu)$  is a convention/scheme choice; only RG-invariant combinations (e.g. transmutation scales) are candidates for physical statements.

**Heuristic PA-H2.11a (Guardrail: geometric weight  $\neq$  anomalous dimension).** The half-density square-root Jacobian is a geometric transformation law under coordinate changes; anomalous dimensions are RG scaling data in interacting theories. They should not be conflated: allowing  $\sigma_*(\mu)$  to run is a representation convenience, not a claim that geometry “produces” anomalous scaling.

## 4. Stationary Phase Produces Half-Density Prefactors (Short-Time Kernel)

In the standard path-integral formalism, stationary phase explains why classical extremals dominate refinement limits. Here we add the complementary kernel-level fact: stationary phase does not only pick the extremal; it also produces a determinant prefactor that transforms as a half-density, i.e. the object needed for coordinate-free kernel composition.

**Derivation PA-D1.4 (Van Vleck prefactor is a bi-half-density).** Let  $S_{\text{cl}}(x, z; t)$  be the classical action as a function of endpoints and time, treated as a generating function. The standard short-time/stationary-phase approximation to the propagator has the form

$$K(x, z; t) \approx \frac{1}{(2\pi i \hbar)^{d/2}} \left| \det \left( -\frac{\partial^2 S_{\text{cl}}}{\partial x \partial z} \right) \right|^{1/2} \exp \left( \frac{i}{\hbar} S_{\text{cl}}(x, z; t) \right).$$

Under a change of coordinates  $x = x(x')$ ,  $z = z(z')$ , the mixed Hessian transforms by the chain rule, and its determinant acquires Jacobian factors:

$$\det\left(-\frac{\partial^2 S_{\text{cl}}}{\partial x' \partial z'}\right) = \det\left(\frac{\partial x}{\partial x'}\right) \det\left(\frac{\partial z}{\partial z'}\right) \det\left(-\frac{\partial^2 S_{\text{cl}}}{\partial x \partial z}\right).$$

Taking square roots shows that the prefactor transforms with  $|\det(\partial x/\partial x')|^{1/2} |\det(\partial z/\partial z')|^{1/2}$ , i.e. exactly as a half-density factor at each endpoint. Thus the stationary-phase prefactor is naturally interpreted as making  $K$  a half-density in each variable, so that kernel composition does not depend on a background measure choice. This is the standard “Van Vleck type” semiclassical prefactor in the correspondence/semiclassical tradition [VanVleck1928Correspondence].

**Derivation PA-D1.4a (Free-propagator semigroup: the  $d/2$  exponent is forced by composition).** The free quantum propagator on  $\mathbb{R}^d$ ,  $K(x, z; t) = \left(\frac{m}{2\pi i \hbar t}\right)^{d/2} \exp\left(\frac{im|x-z|^2}{2\hbar t}\right)$ , provides a concrete witness. As a bi-half-density, the product  $\mathbf{K}(x, y; t_1) \mathbf{K}(y, z; t_2)$  pairs the two  $|dy|^{1/2}$  factors into a density  $|dy|$ , which integrates without a background measure. The  $d$ -dimensional Gaussian integral over  $y$  gives a volume factor  $\left(\frac{2\pi i \hbar t_1 t_2}{m(t_1 + t_2)}\right)^{d/2}$ . Combining with the two kernel prefactors:

$$\left(\frac{m}{2\pi i \hbar t_1}\right)^{d/2} \left(\frac{m}{2\pi i \hbar t_2}\right)^{d/2} \left(\frac{2\pi i \hbar t_1 t_2}{m(t_1 + t_2)}\right)^{d/2} = \left(\frac{m}{2\pi i \hbar (t_1 + t_2)}\right)^{d/2},$$

which is the prefactor of  $K(x, z; t_1 + t_2)$ . The semigroup property holds because the  $d/2$  exponent appears three times (twice from the kernels, once from the Gaussian) and the cancellation is exact only for this exponent. This is the Van Vleck determinant in disguise: for the free particle,  $\det(-\partial^2 S_{\text{cl}}/\partial x \partial z) = (m/t)^d$ , so the square root is  $(m/t)^{d/2}$ .

In the tangent-groupoid near-diagonal picture, the parameter  $\varepsilon = \hbar t/m$  (diffusion scale) plays the role of the rescaling parameter in PA-D1.2a: the prefactor is proportional to  $\varepsilon^{-d/2}$ , and the passage from kernel to symbol on  $TM$  absorbs this half-density Jacobian.

**Derivation PA-D1.9 (Square-root delta normalization has half-density weight).** In finite dimension, the “localize on critical points” distribution is  $\delta(\nabla f)$ , supported on  $\text{Crit}(f)$ . A concrete way it appears is via a “halved” oscillatory integral with a normalization exponent fixed by dimension.

Let  $f : \mathbb{R}^N \rightarrow \mathbb{R}$  be smooth and define, for  $\varepsilon > 0$ ,

$$A_\varepsilon(O) := \varepsilon^{-N/2} \int_{\mathbb{R}^N} e^{\frac{i}{\varepsilon} f(x)} O(x) dx.$$

Then

$$|A_\varepsilon(O)|^2 = \varepsilon^{-N} \iint e^{\frac{i}{\varepsilon} (f(x) - f(y))} O(x) \overline{O(y)} dx dy.$$

Applying the near-diagonal scaling  $y = x + \varepsilon z$  (so  $dy = \varepsilon^N dz$ ) gives

$$|A_\varepsilon(O)|^2 = \iint e^{-\frac{i}{\varepsilon}(f(x+\varepsilon z)-f(x))} O(x) \overline{O(x+\varepsilon z)} dz dx.$$

Formally letting  $\varepsilon \rightarrow 0$  yields

$$|A_\varepsilon(O)|^2 \rightarrow \iint e^{-iz \cdot \nabla f(x)} |O(x)|^2 dz dx = (2\pi)^N \int \delta(\nabla f(x)) |O(x)|^2 dx.$$

The exponent  $N/2$  in the prefactor is exactly the half-density scaling: it cancels the Jacobian  $dy = \varepsilon^N dz$  under near-diagonal rescaling, and it is the “square root” of the density normalization that produces  $\delta(\nabla f)$ .

**Heuristic PA-H1.4 (Where Planck area can enter, minimally).** Derivation PA-D1.3 isolates one minimal route by which a Planck-scale quantity can enter: if the theory supplies a single universal coupling with dimension of length (Newton’s constant) and one demands that the half-density normalization constant be built from that coupling *without fractional powers*, then  $d = 4$  is singled out and the resulting constant has the dimension of an area, naturally identified with the Planck area  $L_P^2 \sim \hbar G_4/c^3$ .

## 5. A Gravitational Anchor: Minimal Areal Speed and the $D = 4$ Cancellation

Rivero’s “Planck areal speed” observation gives a concrete route by which Planck-scale discreteness reappears at Compton scales in inverse-square gravity [RiveroAreal] [RiveroSimple].

**Heuristic PA-H1.3 (Areal-speed selection).** In  $3 + 1$  Newtonian gravity (inverse-square), imposing a discrete areal-speed/area-time condition at a Planck scale can yield characteristic radii proportional to a reduced Compton length, with Newton’s constant canceling when expressed in Planck units. This is a nontrivial indication that “a universal area scale” can be operationally meaningful at low energies in  $D = 4$ .

**Derivation PA-D1.5 (Inverse-square circular orbit + Planck areal speed  $\rightarrow$  Compton radius).** For a circular orbit under an inverse-square central force  $F(r) = K/r^2$  (with coupling  $K > 0$ ), the centripetal balance is  $mv^2/r = K/r^2$ . The areal speed is  $\dot{A} = \frac{1}{2}rv$ , so  $v = 2\dot{A}/r$ . Substituting into the force balance gives

$$m \left( \frac{2\dot{A}}{r} \right)^2 = \frac{K}{r} \implies r = \frac{4m\dot{A}^2}{K}.$$

For Newtonian gravity between a source mass  $M$  and test mass  $m$ ,  $K = GMm$ , hence



$$r = \frac{4\dot{A}^2}{GM},$$

independent of the test mass  $m$ . If one now imposes  $\dot{A} = k \dot{A}_P$ , where Rivero's Planck areal speed is  $\dot{A}_P = cL_P$  [RiveroAreal], then using  $L_P^2 = G\hbar/c^3$  yields

$$r = \frac{4k^2(cL_P)^2}{GM} = \frac{4k^2(G\hbar/c)}{GM} = 4k^2 \frac{\hbar}{cM}.$$

Thus  $r$  becomes a multiple of the reduced Compton length  $L_M = \hbar/(cM)$ , with Newton's constant canceled out. In particular,  $k = \frac{1}{2}$  gives  $r = L_M$ . This is the “Planck area per Planck time  $\Rightarrow$  Compton scale” cancellation highlighted in [RiveroAreal] and summarized in [RiveroSimple].

**Remark PA-D1.5a (Generic  $\backslash(F=K/r^q\backslash)$ : only  $\backslash(q=2\backslash)$  yields linear Compton scaling;  $\backslash(q\backslash)$  links to dimension).** For a power-law central force  $F(r) = K/r^q$  with  $K > 0$ , circular balance gives  $mv^2/r = K/r^q$ , i.e.  $mv^2 = Kr^{1-q}$ . Using the circular areal speed  $\dot{A} = \frac{1}{2}rv$  (so  $v = 2\dot{A}/r$ ) yields

$$4m\dot{A}^2 = Kr^{3-q}.$$

Hence, for  $q \neq 3$ ,

$$r = \left( \frac{4m\dot{A}^2}{K} \right)^{\frac{1}{3-q}},$$

while for  $q = 3$  the radius drops out and  $4m\dot{A}^2 = K$ .

In the gravitational specialization  $K = GMm$ , the test mass cancels as before and

$$r = \left( \frac{4\dot{A}^2}{GM} \right)^{\frac{1}{3-q}}.$$

If one imposes  $\dot{A} = cL_P$  and uses the  $D = 4$  identity  $L_P^2 = G\hbar/c^3$ , then

$$r = \left( \frac{4\hbar}{cM} \right)^{\frac{1}{3-q}}.$$

Thus the Planck-areal-speed substitution produces *linear* reduced-Compton scaling  $r \propto \hbar/(cM)$  only for  $q = 2$  (inverse-square). For Newtonian long-range fields in  $n$  spatial dimensions, the Laplacian Green function gives  $\Phi(r) \propto r^{2-n}$  (for  $n > 2$ ), so  $F \sim |\nabla\Phi| \propto r^{1-n}$ , i.e.  $q = n - 1$  (with the  $n = 2$  logarithmic exception) [Tanaka2021KernelQuadrature]. In this sense the  $q = 2$  special case corresponds to  $n = 3$  spatial dimensions (spacetime  $D = 4$ ). Equivalently, substituting  $q = n - 1$  into the mass scaling gives  $r \propto M^{-1/(4-n)}$  (for  $n \neq 4$ ), so the linear Compton scaling (and the  $G$ -cancellation in the  $D = 4$  identity  $L_P^2 = G\hbar/c^3$ ) is uniquely  $n = 3$ ; the  $n = 4$  case is the degenerate  $q = 3$  condition where the radius drops out.

Remark PA-D1.5b (SR continuation of the inverse-square witness: Compton branch  $\backslash(\backslash\text{to}\backslash)$  Planck floor). Inside a mechanical special-relativistic model with an external inverse-square force (no GR field dynamics), one can keep the fixed coordinate-time areal speed  $\dot{A}_0 = dA/dt$  and continue PA-D1.5 exactly.

For inverse-square forces, SR circular motion gives  $v = K/L$ , while

$$L = \gamma m r v = 2\gamma m \dot{A}_0, \quad \gamma = \frac{1}{\sqrt{1 - v^2/c^2}}.$$

Hence

$$\gamma v = \frac{K}{2m\dot{A}_0},$$

so

$$v = \frac{\frac{K}{2m\dot{A}_0}}{\sqrt{1 + \left(\frac{K}{2m\dot{A}_0 c}\right)^2}}, \quad r = \frac{2\dot{A}_0}{v} = \frac{4m\dot{A}_0^2}{K} \sqrt{1 + \left(\frac{K}{2m\dot{A}_0 c}\right)^2}.$$

For gravity  $K = GMm$ ,  $m$  cancels:

$$r(M, \dot{A}_0) = \frac{4\dot{A}_0^2}{GM} \sqrt{1 + \left(\frac{GM}{2\dot{A}_0 c}\right)^2}.$$

With  $\dot{A}_0 = cL_P$ , using  $L_P^2 = \hbar G/c^3$  and  $M_P^2 = \hbar c/G$ ,

$$r(M) = \frac{4\hbar}{Mc} \sqrt{1 + \frac{M^2}{4M_P^2}}.$$

Therefore  $M \ll M_P$  reproduces the PA-D1.5 Compton-like branch  $r \approx 4\hbar/(Mc)$ , while  $M \gg M_P$  saturates at

$$r \rightarrow 2L_P.$$

So in this SR continuation, the inverse-square Planck-areal-speed witness is not destroyed; it is regularized into a bounded interpolation.

Remark PA-D1.5c (Clock-choice sensitivity: fixed  $\backslash(dA/dt\backslash)$  vs fixed  $\backslash(dA/d\tau\backslash)$ ). The  $2L_P$  high-mass saturation in PA-D1.5b is tied to fixing the coordinate-time areal speed  $\dot{A}_t = dA/dt$ . If one instead fixes proper-time areal speed  $\dot{A}_\tau = dA/d\tau$  in the same inverse-square SR model, then

$$r(M, \dot{A}_\tau) = \frac{4\dot{A}_\tau^2}{GM} \sqrt{1 - \left(\frac{GM}{2\dot{A}_\tau c}\right)^2},$$

which is defined only for  $GM < 2\dot{A}_\tau c$ . With  $\dot{A}_\tau = cL_P$ ,

$$r(M) = \frac{4\hbar}{Mc} \sqrt{1 - \frac{M^2}{4M_P^2}},$$

so  $M \rightarrow 2M_P^- \Rightarrow r \rightarrow 0$ , not  $2L_P$ . Therefore the low-mass Compton-like branch is robust, while the high-mass asymptotic is clock-convention dependent in this mechanical SR setting.

**Remark PA-D1.5d (Invariant-candidate reformulation via specific angular momentum).** A clock-independent candidate in the same central-source setup is the specific angular momentum scalar

$$\ell := \frac{1}{m} \sqrt{\frac{1}{2} L_{\mu\nu} L^{\mu\nu}},$$

where  $L^{\mu\nu}$  is the source-rest-space projection of orbital  $J^{\mu\nu}$  (using source 4-velocity  $U^\mu$ ). In the source rest frame this reduces to

$$\ell = \gamma r v, \quad \dot{A}_t = \frac{\ell}{2\gamma}, \quad \dot{A}_r = \frac{\ell}{2}.$$

So fixing  $\ell$  selects the proper-time branch rather than the coordinate-time branch. For inverse-square gravity,

$$r(M, \ell) = \frac{\ell^2}{GM} \sqrt{1 - \left(\frac{GM}{\ell c}\right)^2}, \quad \ell > \frac{GM}{c},$$

and  $\ell = 2cL_P$  reproduces PA-D1.5c. This does not yet prove universality, but it provides a structurally covariant way to encode the postulate without choosing a clock variable directly.

**Remark PA-D1.5e (Why simple frame-free bivector invariants are insufficient here).** One might try to avoid the source-velocity projection entirely and use only Lorentz invariants of

$$M^{\mu\nu} = R^\mu p^\nu - R^\nu p^\mu,$$

namely

$$I_1 = \frac{1}{2} M_{\mu\nu} M^{\mu\nu} = R^2 p^2 - (R \cdot p)^2, \quad I_2 = \frac{1}{2} M_{\mu\nu} \star M^{\mu\nu}.$$

But for the circular central branch ( $R \cdot p = 0$ ,  $R^2 = r^2$ ,  $p^2 = -m^2 c^2$ ), this gives

$$I_1 = -m^2 c^2 r^2,$$

independent of orbital speed, while  $I_2$  vanishes in the planar case. So these simple frame-free invariants do not encode the areal-rate branch parameter. In this setup, a timelike direction (e.g. source  $U^\mu$ ) appears to be minimal extra structure for a useful covariant postulate.

**Remark PA-D1.5f (Minimal timelike-structure rule for this branch).** For the present central-source inverse-square model, a practical “minimal structure” rule is: 1. use the source worldline 4-velocity  $U^\mu$  as the distinguished

timelike direction (already part of the model input), and 2. formulate the postulate on

$$\ell_{(U)} := \frac{1}{m} \sqrt{\frac{1}{2} M_{\mu\nu}^{(U)} M_{(U)}^{\mu\nu}}, \quad M^{(U)\mu\nu} = h(U)^\mu{}_\alpha h(U)^\nu{}_\beta M^{\alpha\beta},$$

rather than directly on coordinate-time areal rate. In the source rest frame this is equivalent to fixing  $\dot{A}_{\tau,U} = \ell_{(U)}/2$ , while coordinate-time rates are derived via the corresponding lapse factor. This keeps the branch covariant-with-source and avoids introducing an additional arbitrary observer field  $u^\mu(x)$ .

**Remark PA-D1.5g (Non-circular planar extension is kinematic).** The  $\ell$ -based rule is not restricted to circular trajectories. For general planar motion with tangential component  $v_\perp = r\dot{\phi}$ ,

$$\ell = \gamma r v_\perp = \gamma r^2 \dot{\phi}, \quad \frac{dA}{dt} = \frac{1}{2} r^2 \dot{\phi} = \frac{\ell}{2\gamma}, \quad \frac{dA}{d\tau} = \frac{\ell}{2}.$$

Thus “fix  $\ell$ ” remains equivalent to fixing proper-time areal rate in the source frame even away from circular orbits; this part is kinematic and does not depend on the specific force law.

**Remark PA-D1.5h (Non-planar caution: vector area rate is fundamental).** For general 3D motion the natural identity is vector-valued:

$$\boldsymbol{\ell} = \frac{1}{m} (\mathbf{r} \times \mathbf{p}) = \gamma (\mathbf{r} \times \mathbf{v}), \quad \frac{d\mathbf{A}}{d\tau} = \frac{\boldsymbol{\ell}}{2}.$$

So the  $\ell$ -rule still survives kinematically, but scalar areal rates require a chosen normal  $\mathbf{n}$ :

$$\frac{dA_{\mathbf{n}}}{d\tau} = \frac{\boldsymbol{\ell} \cdot \mathbf{n}}{2}.$$

In non-planar perturbations, interpreting a scalar “areal speed postulate” without specifying this projection is ambiguous; the projection choice is part of the model specification.

**Remark PA-D1.5i (Observability criterion for projected areal-rate postulates).** In perturbed-orbit settings, a projected areal-rate claim is empirically meaningful only after specifying: 1. projection normal  $\mathbf{n}$ , 2. clock convention ( $t$ -based or  $\tau$ -based), 3. reconstruction map for  $(\mathbf{r}, \mathbf{v})$  in the source frame. With those choices fixed, the observable is

$$\dot{A}_{\mathbf{n}}(t) = \frac{1}{2} \mathbf{n} \cdot (\mathbf{r} \times \mathbf{v}), \quad \frac{dA_{\mathbf{n}}}{d\tau} = \frac{1}{2} \mathbf{n} \cdot \boldsymbol{\ell},$$

and departures from constant projected areal rate are governed by projected torque

$$\frac{d}{dt} (\mathbf{n} \cdot \boldsymbol{\ell}) = \frac{1}{m} \mathbf{n} \cdot (\mathbf{r} \times \mathbf{F}).$$

So the postulate becomes falsifiable precisely when projection, clock, and reconstruction are part of the model declaration.

**Remark PA-D1.5j (Minimal implementation pipeline).** A compact data-to-test pipeline is: 1. reconstruct object state in observer frame from direction  $\hat{\mathbf{n}}(t)$ , range  $\rho(t)$ , and line-of-sight velocity  $\dot{\rho}(t)$ :

$$\mathbf{r}_{\text{obj}} = \rho \hat{\mathbf{n}}, \quad \mathbf{v}_{\text{obj}} = \dot{\rho} \hat{\mathbf{n}} + \rho \dot{\hat{\mathbf{n}}},$$

2. subtract source ephemeris to obtain source-frame relative state  $(\mathbf{r}, \mathbf{v})$ , 3. evaluate

$$\dot{A}_{\mathbf{n}} = \frac{1}{2} \mathbf{n} \cdot (\mathbf{r} \times \mathbf{v}), \quad \frac{dA_{\mathbf{n}}}{d\tau} = \frac{1}{2} \mathbf{n} \cdot \boldsymbol{\ell},$$

and, if a force model is supplied, the projected-torque residual

$$\mathcal{T}_{\mathbf{n}} := \frac{d}{dt}(\mathbf{n} \cdot \boldsymbol{\ell}) - \frac{1}{m} \mathbf{n} \cdot (\mathbf{r} \times \mathbf{F}).$$

This keeps the postulate test tied to explicit reconstruction and uncertainty handling rather than to abstract kinematic statements alone.

**Remark PA-D1.5k (Minimal uncertainty scaffold).** At first order, uncertainty in projected observables can be propagated by Jacobians:

$$q := \dot{A}_{\mathbf{n}} \Rightarrow \sigma_q^2 \approx J_q \Sigma_x J_q^\top,$$

for reconstructed state vector  $x$  and covariance  $\Sigma_x$ , and

$$\sigma_{\mathcal{T}}^2 \approx J_{\mathcal{T}} \Sigma_z J_{\mathcal{T}}^\top$$

for residual  $\mathcal{T}_{\mathbf{n}}$  with augmented state  $z$ . This linear scaffold is a baseline; in strongly nonlinear regimes the same quantities should be cross-checked with nonlinear propagation (e.g. Monte Carlo) before interpretation.

**Remark PA-D1.5l (Practical nonlinear-validation trigger).** A lightweight policy is to run a pilot nonlinear propagation check and compare against the linear  $\sigma$  estimate; if the discrepancy is at the few-percent level (or larger), treat linearized errors as insufficient and switch to nonlinear uncertainty propagation for reporting.

**Remark PA-D1.5m (Regime-dependent trigger calibration).** The trigger in PA-D1.5l should be calibrated by uncertainty regime, not treated as universal. A practical diagnostic pair is

$$\epsilon_{\text{nl}} := \frac{|\sigma_{\text{MC}} - \sigma_{\text{lin}}|}{\sigma_{\text{lin}}}, \quad \chi := \max\left(\frac{\sqrt{\text{tr } \Sigma_r}}{\|\mathbf{r}\|}, \frac{\sqrt{\text{tr } \Sigma_v}}{\|\mathbf{v}\|}\right),$$

with  $\Sigma_r, \Sigma_v$  the position/velocity covariance blocks in the chosen reconstruction model. Pilot scans in correlated-noise families can then map  $\epsilon_{\text{nl}}(\chi)$  for the instrument/model pair; the “few-percent” policy corresponds to selecting an operational  $\epsilon_{\text{nl}}$  band after this calibration, rather than imposing a context-free constant.

## 6. Connection to the Refinement-Composition Framework

The broader program in which this note sits argues that: 1. classical dynamics are recovered from quantum composition by stationary-phase concentration, and 2. refinement across scales forces RG-style consistency conditions when naive limits diverge.

This note adds a complementary ingredient: the kernel side is most naturally formulated in half-density language, and stationary phase produces the bi-half-density prefactor directly. A universal convention for turning those half-densities into scalar amplitudes then requires a length <sup>$d/2$</sup>  scale; in  $d = 4$  this is an area scale.

## 7. Open Problems and Outlook

1. (*Addressed: Derivation PA-D1.4a.*) The free-propagator semigroup computation provides a concrete witness; a full tangent-groupoid treatment remains desirable.
2. (*Addressed: Remark PA-E5c.*) The Van Vleck / transmutation separation is explicit in the 2D delta model; a  $d = 4$  gauge-theory witness is still open.
3. General-dimension analysis: clarify what replaces “area” in odd dimensions and whether a universal normalization is still defensible. (*Partially addressed: Remark PA-D1.3a classifies the three obstruction mechanisms. An explicit odd-d transmutation witness is still open.*)
4. (*Addressed: Remark PA-H2.6b.*) The minimal hypothesis set is PA-H2.1 + PA-H2.4 + PA-H2.5a + gravity-only; each is necessary.
5. Track minimal-length/GUP scenarios as a comparison branch: do they implement the “needed scale” at the level of kinematics (modified commutators/dispersion) or can they be reframed as a refinement-compatibility condition? Use [Hossenfelder2013MinimalLength] as an OA entry point.

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