

Scale-Invariant Information Limit in Nuclear Optical Potential Extraction

Jin Lei^{1,*}

¹*School of Physics Science and Engineering, Tongji University, Shanghai 200092, China*

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How strongly can nucleon-nucleus scattering data constrain optical model potential parameters? I show that the continuous “Igo ambiguity” reflects a physical information limit, not insufficient data. Using Fisher information geometry with the 11-parameter Koning-Delaroche potential (including spin-orbit), I decompose the constraining power step by step: single-energy elastic scattering constrains only $D_{\text{eff}} \approx 1.7 \pm 0.4$ parameter combinations averaged over 168 configurations; at a given energy, adding the reaction cross section σ_R provides negligible improvement for comparable fractional uncertainties (a single datum versus 17 angular bins), while the analyzing power A_y lifts D_{eff} by $\sim 30\%$ through spin-orbit information; combining measurements across 7 energies provides the largest improvement, raising D_{eff} to ~ 2.2 . Crucially, this limit is *scale-invariant*: under uniform error reduction, better experiments shrink error bars but cannot resolve the underlying parameter degeneracy, just as shorter-wavelength light improves resolution but cannot beat the diffraction limit for a given aperture. Verified across 12 nuclei ($A = 12\text{--}208$), 7 energies ($E = 10\text{--}200$ MeV), and both projectile types, these results provide the first quantitative characterization of the Igo ambiguity and establish a clear hierarchy of experimental measurements for constraining the optical potential.

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Introduction.—The extraction of fundamental interactions from scattering observables represents one of the quintessential inverse problems in quantum physics. Since the inception of the optical model potential (OMP) over seventy years ago [1–3], this effective description has served as the cornerstone for nuclear reaction theory. Yet, despite the exponential growth in experimental precision and computational power, a fundamental puzzle persists: the “inverse problem ill-posedness,” historically known as the Igo ambiguity [4]. Distinct parameter sets, representing vastly different nuclear potentials, can yield indistinguishable scattering cross sections, creating a continuous degeneracy that defies unique determination.

Two types of ambiguity exist [5]: *discrete* ambiguity from different wavefunction node numbers (resolvable by high-energy rainbow scattering), and *continuous* ambiguity where parameters trade off while maintaining the surface potential value. The qualitative origin of this continuous ambiguity is well understood: elastic scattering is primarily sensitive to the potential near the nuclear surface, making it insensitive to the detailed radial shape [5]. However, three key questions remain unanswered: (1) *How many* independent parameter combinations can actually be constrained? (2) Can improved experimental precision eventually resolve the degeneracy? (3) Is the ambiguity universal across the nuclear chart?

Recent Bayesian uncertainty quantification has documented large parameter uncertainties. In a landmark comparison, King *et al.* [6] showed that Bayesian MCMC posterior distributions for optical potential parameters are strongly non-Gaussian, with only the depth V and radius r_v strongly correlated—in striking contrast to the frequentist χ^2 approach which produced spurious correlations among all parameters. Subsequent studies con-

firmed these findings [7, 8], while principal component analysis (PCA) identified the dominant variance directions [9]. Yet these empirical observations, while valuable, require $\sim 100,000$ MCMC samples per system and lack a rigorous theoretical framework that would answer the questions above.

In this Letter, I provide the first quantitative characterization of the Igo ambiguity using Fisher information geometry. By progressively adding observables and energies, I decompose the constraining power of nucleon-nucleus scattering data, revealing that elastic scattering compresses the 11-dimensional parameter space into a drastically lower-dimensional subspace, with only ~ 2 “stiff” parameter combinations constrainable while the remaining directions remain “sloppy.”

Using Fisher information geometry, I demonstrate that single-energy elastic scattering universally constrains only

$$D_{\text{eff}} \approx 1.7 \pm 0.4 \quad (1)$$

independent parameter combinations out of 11 in the Koning-Delaroche potential [10] (including spin-orbit coupling). This result holds regardless of target nucleus ($A = 12\text{--}208$) or beam energy ($E = 10\text{--}200$ MeV). Crucially, this limit is *scale-invariant*: improving experimental precision reduces error bars but cannot alter the “flat” topology of the parameter manifold. Just as increasing the pixel count cannot resolve features below the diffraction limit, reducing experimental errors cannot resolve parameter degeneracy beyond this information bound.

All calculations employ the Numerov algorithm for solving the radial Schrödinger equation, with finite-difference gradients validated against the coupled-channels code FRESKO [11]. The analysis spans both

spin-0 (9-parameter central potential) and spin-1/2 (11-parameter with spin-orbit) scattering, incorporating elastic cross sections, analyzing powers, and reaction and total cross sections.

Method.—Consider nucleon-nucleus elastic scattering with the Koning-Delaroche (KD02) optical potential [10], which has 9 parameters for the central terms: real volume (V, r_v, a_v), imaginary volume (W, r_w, a_w), and imaginary surface (W_d, r_d, a_d). Including spin-orbit coupling with Thomas form adds 2 parameters (V_{so}, W_{so}) for the real and imaginary spin-orbit depths, with the spin-orbit geometry (r_{so}, a_{so}) fixed at the KD02 values. This yields an 11-parameter model for spin-1/2 scattering.

The question is: how much can elastic cross section data $\sigma(\theta)$ tell us about these 11 parameters? To answer this, I compute the *relative sensitivity* of the cross section to each parameter:

$$S_i(\theta) = \frac{\partial \log \sigma(\theta)}{\partial \log p_i} = \frac{p_i}{\sigma} \frac{\partial \sigma}{\partial p_i}. \quad (2)$$

Using logarithmic derivatives makes parameters with different physical units (e.g., $V \sim 50$ MeV vs $r_v \sim 1.2$ fm) directly comparable. If S_V and S_{r_v} point in nearly the same direction in function space, then a fractional increase in V has nearly the same effect as a fractional increase in r_v , and the two cannot be independently determined.

This is formalized through the Fisher information matrix (FIM):

$$F_{ij} = \sum_{\theta} S_i(\theta) S_j(\theta), \quad (3)$$

where the sum runs over angular bins of the elastic cross section, assuming uniform relative uncertainties. More generally, for a data vector $\{y_k\}$ comprising $d\sigma/d\Omega$, A_y , σ_R , and σ_T , the FIM takes the standard form $F_{ij} = \sum_k \sigma_k^{-2} (\partial y_k / \partial p_i) (\partial y_k / \partial p_j)$. For positive-definite cross sections ($d\sigma/d\Omega$, σ_R , σ_T), relative uncertainties recover the log-derivative form; for A_y (which can vanish or change sign), absolute uncertainties are used ($\Delta A_y = 0.03$, a typical experimental scale). Each observable type enters additively ($F = F_{\text{elastic}} + F_{A_y} + \dots$), enabling the step-by-step decomposition below. This convention makes D_{eff} a geometric property of the scattering problem rather than experimental details. The FIM encodes how sensitively observables respond to parameter changes. Diagonalizing the FIM yields eigenvalues $\{\lambda_i\}$ and corresponding eigenvectors $\{\mathbf{e}_i\}$. Each eigenvalue measures the information content along its eigenvector direction, with the fractional information $\lambda_i / \sum_j \lambda_j$. Each eigenvector $\mathbf{e}_i = (e_{i,1}, \dots, e_{i,N_p})$ defines a linear combination of parameters; the squared component $e_{i,k}^2$ gives the fractional contribution of parameter k to that mode. Large eigenvalues indicate “stiff” directions (well-constrained), while small eigenvalues indicate “sloppy” directions (poorly constrained).

The effective dimensionality counts how many stiff directions exist:

$$D_{\text{eff}} = \frac{(\sum_i \lambda_i)^2}{\sum_i \lambda_i^2}. \quad (4)$$

This equals N_p (the number of parameters) if all directions are equally stiff, and approaches 1 if one direction dominates. Physically, D_{eff} answers: “How many independent parameter combinations can the data actually constrain?”

Geometrically, the FIM represents the local curvature of the χ^2 landscape: large eigenvalues correspond to “stiff” directions (well-constrained), small eigenvalues to “sloppy” directions (the Igo ambiguity). Finding $D_{\text{eff}} \approx 1.7$ means the 11-dimensional parameter space collapses into a ~ 9 -dimensional degenerate manifold.

Crucially, D_{eff} is *scale-invariant* under uniform rescaling of experimental uncertainties. Reducing experimental error ϵ scales all eigenvalues as $\lambda_i \propto 1/\epsilon^2$, which tightens all error bars proportionally. But the *ratio* of eigenvalues, and hence D_{eff} , is unchanged. Better precision shrinks the allowed parameter region but cannot change its shape from a narrow valley into a sphere.

Numerical method.—The Schrödinger equation for each partial wave is solved using the Numerov algorithm with matching to Coulomb functions at large radius. For spin-1/2 scattering, the solver iterates over all (l, j) partial wave channels, with the spin-orbit potential entering through the Thomas form with factor of $2(\mathbf{l} \cdot \boldsymbol{\sigma} = 2\mathbf{l} \cdot \mathbf{s})$. From the (l, j) -resolved S -matrix, the direct and spin-flip scattering amplitudes $f(\theta)$ and $g(\theta)$ are constructed, yielding the differential cross section $d\sigma/d\Omega = |f|^2 + |g|^2$, analyzing power $A_y = 2\text{Im}(fg^*)/(|f|^2 + |g|^2)$, reaction cross section σ_R , and total cross section σ_T .

Parameter sensitivities are computed using central finite differences: $\partial \sigma / \partial p_i \approx [\sigma(p_i + \delta) - \sigma(p_i - \delta)] / (2\delta)$ with $\delta = 0.01|p_i|$. For the leading eigenvalues that determine D_{eff} , this relative perturbation provides sufficient numerical accuracy; the D_{eff} values are robust to order-of-magnitude variations in δ . The Numerov solver has been validated against the coupled-channels code FRESKO [11], achieving S -matrix agreement to $|\Delta S| < 0.002$ and cross-section agreement to $< 0.2\%$.

Results.—I systematically analyze 168 configurations: 12 nuclei (^{12}C to ^{208}Pb), 7 energies (10–200 MeV), and both neutron and proton projectiles. Figure 1 displays D_{eff} as a heatmap across all configurations. The remarkable uniformity ($D_{\text{eff}} \approx 1\text{--}3$ everywhere) demonstrates that the information limit is universal: it depends neither on target mass nor beam energy nor projectile type. The statistics yield $D_{\text{eff}} = 1.63 \pm 0.34$ (neutrons) and 1.73 ± 0.52 (protons).

Figure 2 confirms this universality through systematic analysis. Panel (a) shows no significant mass dependence from ^{12}C to ^{208}Pb at $E = 50$ MeV. Panel (b) shows D_{eff} remains bounded between 1 and 3 across all energies for

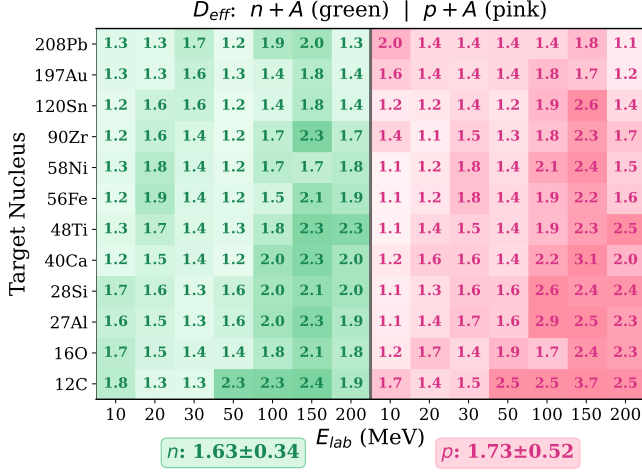


FIG. 1. Universal information limit. Effective dimensionality D_{eff} across 12 nuclei and 7 energies for neutron (green) and proton (pink) projectiles. Values cluster around 1–3 regardless of system, demonstrating that the information limit is intrinsic to elastic scattering.

$n+^{40}\text{Ca}$. Panel (c) reveals condition numbers (ratio of largest to smallest FIM eigenvalue) exceeding 10^7 , indicating an extreme hierarchy between constrained and unconstrained directions. This means the stiffest direction is ten million times more constrained than the sloppiest: constraining “stiff” parameters to 1% precision requires standard experimental errors, but constraining “sloppy” parameters to the same level would require reducing errors by a factor of $\sqrt{10^7} \approx 3000$. For all practical purposes, the ~ 9 “sloppy” dimensions are unobservable.

Anatomy of the information limit.—What physics do these ~ 2 constrained directions represent? Figure 3 shows the eigenvector composition for $n+^{40}\text{Ca}$ at 50 MeV using the 11-parameter model with elastic angular distributions. The dominant eigenvector \mathbf{e}_1 (92% of information) is composed primarily of r_v (65%), r_d (15%), and V (12%). Crucially, V and r_v enter \mathbf{e}_1 with the *same sign*. This identifies \mathbf{e}_1 as the “volume integral” direction: V and r_v both increasing together makes $J_V \propto V r_v^3$ larger, producing a measurably different cross section. In contrast, the Igo ambiguity [4] lives in the *orthogonal* sloppy direction, where V increases while r_v decreases so that J_V stays approximately constant. This sloppy direction has an eigenvalue 10^5 times smaller—the data cannot distinguish a deep, narrow potential from a shallow, wide one if both produce the same scattering surface. Stated simply: there are two directions in the V - r_v plane; the data pin down the volume integral (stiff), but cannot resolve the depth-radius tradeoff (sloppy).

The radius r_v dominates \mathbf{e}_1 over V because $J_V \propto V r_v^3$: in logarithmic derivatives, $\partial \log J_V / \partial \log r_v = 3$ versus $\partial \log J_V / \partial \log V = 1$. The second eigenvector \mathbf{e}_2 (6% of information) mixes r_v (29%), a_v (27%), and r_d (27%),

encoding the radial shape of the potential surface. Together, these two modes capture 98% of all information. The remaining 9 directions (diffuseness, imaginary potential, and spin-orbit) contribute only 2% combined—they are effectively “invisible” to elastic angular distributions, confirming that reaction cross-section and polarization measurements are essential for constraining them.

Why do r_v and r_d appear in both eigenvectors? Both the real and imaginary potentials have Woods-Saxon shapes peaked near the nuclear surface. Because the cross section depends on the potential near the classical turning point, the real and surface radii (r_v , r_d) dominate the sensitivity while diffuseness parameters (a_v , a_w , a_d) produce only subtle changes in the radial tail (see End Matter, Fig. 7). The angle-resolved sensitivity curves $|S_V(\theta)|$ and $|S_{r_v}(\theta)|$ have nearly identical angular shapes, differing primarily in magnitude (End Matter, Fig. 6). This geometric collinearity is the origin of the V - r_v degeneracy: because a fractional change in V has nearly the same *angular pattern* as a fractional change in r_v , the data cannot separate their individual contributions. Only the “scattering surface” combination (dominated by r_v due to the cubic power in J_V) is measurable. But can additional observables lift this degeneracy?

Step-by-step constraint analysis.—A key question is whether additional observables and measurements can break the degeneracy. I decompose the constraining power by progressively adding information sources, as shown in Fig. 4 for four representative systems.

Step 1: Elastic $d\sigma/d\Omega$ alone.—Starting from 17 angular bins of the elastic cross section at 50 MeV, $D_{\text{eff}} \approx 1.2$ –1.4 for all systems studied. Only the “potential radius” mode (predominantly r_v) is constrained; the remaining 10 directions are sloppy. The eigenvalue spectrum drops by seven orders of magnitude from the stiffest to the sloppiest direction.

Step 2: Adding σ_R .—The reaction cross section provides one additional datum sensitive to the imaginary potential. However, this single number is overwhelmed by the 17 elastic angular bins: D_{eff} increases by less than 0.001. The information from σ_R is deposited along directions already constrained by elastic scattering, producing no meaningful improvement.

Step 3: Adding A_y and σ_T .—The analyzing power, measured at the same 17 angles, probes spin-orbit interference through the spin-flip amplitude $g(\theta)$. This provides genuinely new information: for $n+^{40}\text{Ca}$, D_{eff} rises from 1.18 to 1.52. The spin-orbit subgroup trace increases by two orders of magnitude, and the information in the imaginary subspace also grows substantially. For $n+^{120}\text{Sn}$, the improvement is even larger ($D_{\text{eff}}: 1.16 \rightarrow 1.66$). The total cross section σ_T provides additional constraint through the optical theorem.

Step 4: Multi-energy combination.—This provides the largest improvement. Combining elastic $d\sigma/d\Omega$ across 7 energies (10–200 MeV) raises D_{eff} from ~ 1.2 to ~ 2.1 –

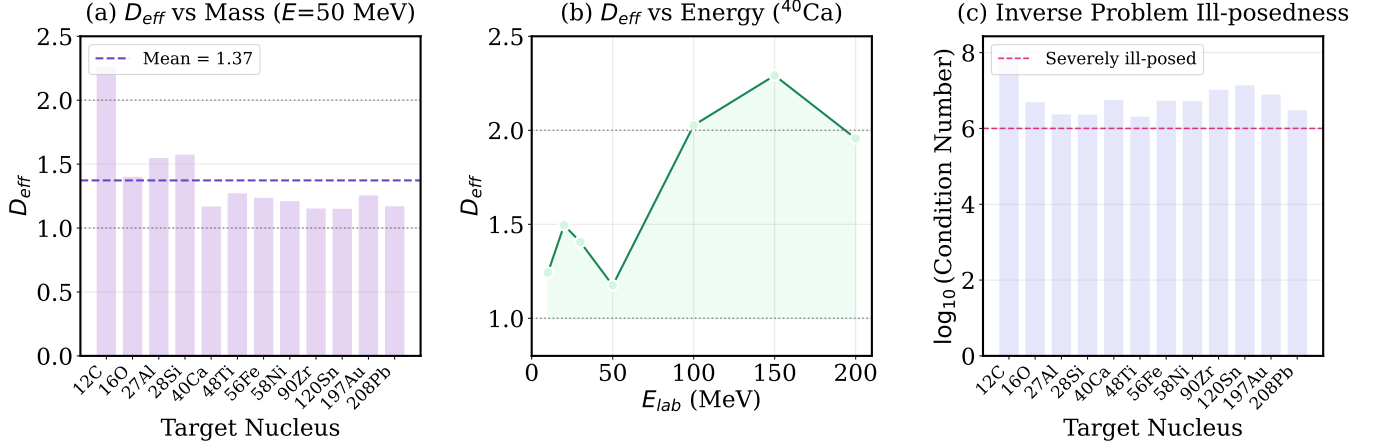


FIG. 2. Systematic analysis for $n + A$ scattering. (a) D_{eff} versus mass at $E = 50$ MeV shows no A -dependence. (b) D_{eff} versus energy for $n + ^{40}\text{Ca}$ stays bounded. (c) Condition number exceeds 10^7 , indicating extreme hierarchy between constrained and unconstrained directions.

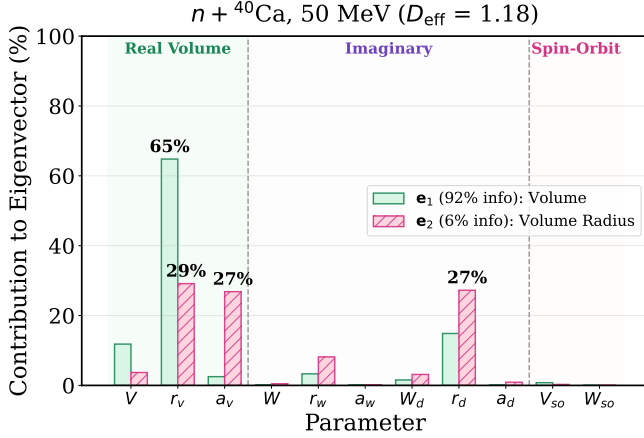


FIG. 3. Eigenvector composition for elastic $n + ^{40}\text{Ca}$ at 50 MeV (11 parameters). The “potential radius” mode \mathbf{e}_1 (green) is dominated by r_v (65%) with r_d (15%) and V (12%); the “radial geometry” mode \mathbf{e}_2 (rose) mixes r_v , a_v , and r_d . Together they capture 98% of the information.

2.3, even without A_y data. The combined Fisher matrix $F_{\text{combined}} = \sum_E F(E)$ accesses the energy dependence of the potential: at low energy the real potential dominates, while at high energy absorption becomes stronger. This energy lever arm lifts the imaginary-parameter degeneracy, with the imaginary subgroup D_{eff} reaching ~ 2.3 (compared to ~ 1.1 at single energy). Adding all observables at all energies pushes D_{eff} to 2.2–2.5, consistent with the success of dispersive optical models [12] that exploit energy-dependent dispersion relations.

Step 5: KD02 global systematics.—Steps 1–4 all constrain the same 11 local parameters at each energy; adding data always improves constraints and raises D_{eff} . Step 5 asks a *different* question. The KD02 global potential [10] expresses each local parameter as a polyno-

mial in energy, e.g., $V(E) = v_1[1 - v_2(E - E_f) + v_3(E - E_f)^2 - v_4(E - E_f)^3]$, introducing a total of 17 global coefficients. Now consider: multi-energy data determine V at 7 energies ($V(10), V(20), \dots, V(200)$ MeV). This is enough to pin down v_1 and v_2 (the overall depth and its linear energy slope), but the cubic polynomial has *four* coefficients. The higher-order terms v_3 and v_4 produce differences that are far smaller than the uncertainties on V itself—the data simply cannot tell a quadratic from a cubic energy dependence. The same redundancy applies to other polynomial chains ($W(E)$, $W_d(E)$, etc.).

The result is $D_{\text{eff}} \approx 1.0$ of 17 global parameters. This does *not* mean that information is lost relative to Step 4—the data still constrain the same local-parameter combinations. Rather, the global KD02 functional form has more adjustable coefficients than the data can resolve, so the additional degrees of freedom simply dilute the effective dimensionality. This has a practical implication: future global optical potentials should use more parsimonious functional forms (fewer higher-order energy terms) to avoid building in degeneracies that the data cannot break.

Angle-resolved sensitivity.—The angular dependence of parameter sensitivity $|S_i(\theta)|$ (End Matter, Fig. 7) reveals that V and r_v have nearly identical angular patterns (explaining the Igo degeneracy), while diffuseness parameters become most distinguishable at backward angles ($\theta > 130^\circ$). The cumulative Fisher information shows that $\sim 90\%$ of the constraining power is captured by $\theta \approx 140^\circ$, providing practical guidance for experimental design: backward-angle measurements are disproportionately valuable for breaking parameter degeneracies.

Discussion.—The step-by-step analysis (Fig. 4) reveals a clear hierarchy of experimental measurements for constraining optical potentials. At 50 MeV, elastic angular distributions alone constrain only $D_{\text{eff}} \approx 1.2$ (the overall

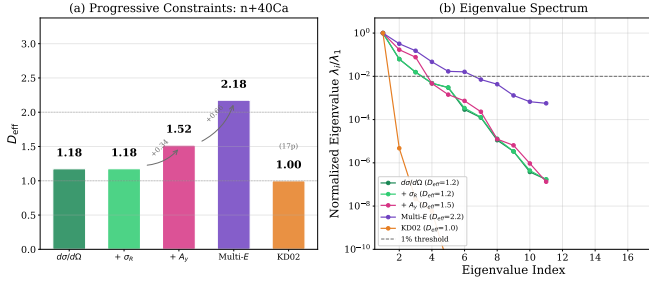


FIG. 4. Step-by-step constraint analysis for $n+^{40}\text{Ca}$. (a) D_{eff} for 11 local parameters increases as observables and energies are added (Steps 1–4): elastic alone (1.18), $+\sigma_R$ (no change), $+A_y$ (1.52), multi-energy (2.18). Step 5 re-expresses the same data in terms of 17 KD02 global parameters: $D_{\text{eff}} \approx 1.0$ of 17 because the polynomial energy dependence has more coefficients than the data can resolve (see text). (b) Eigenvalue spectra at each step.

168-configuration average is 1.7 ± 0.4); σ_R adds negligible information due to being a single integrated quantity; A_y provides genuine new information through spin-orbit interference, lifting D_{eff} by $\sim 30\%$; and multi-energy data provide the largest improvement, reaching $D_{\text{eff}} \approx 2.2$.

The scale-invariance of D_{eff} has important implications. No matter how precisely cross sections are measured, the parameter space retains its “flat valley” topology. This is a geometric property of how scattering encodes potential information, not a limitation of current experiments. Notably, D_{eff} measures how *uniformly* information is distributed across parameter directions, not the total amount; more data always reduce error bars but do not necessarily increase D_{eff} . Indeed, in the 45-dimensional universal KD02 coefficient space, A_y data actually *lower* D_{eff} from 2.5 to 2.0 because they reinforce already-stiff directions rather than opening new ones (see End Matter).

These results provide a rigorous foundation for the empirical findings of King *et al.* [6], who showed using Bayesian MCMC that only V and r_v are strongly correlated. In the Gaussian approximation, the posterior covariance equals the inverse Fisher matrix ($\Sigma_{\text{post}} = F^{-1}$), so $D_{\text{eff}} \approx 1.7$ directly explains why only ~ 2 correlation structures are visible in MCMC scatter plots. A key advantage of the Fisher approach is computational efficiency: D_{eff} requires only $2N_p + 1 = 23$ solver calls versus $\sim 100,000$ MCMC samples, enabling the systematic 168-configuration scan presented here. The detailed Fisher-Bayesian comparison is given in the End Matter.

The step-by-step decomposition provides clear experimental guidance: A_y measurements are more valuable than σ_R at a single energy (17 data points sensitive to spin-orbit versus a single datum), but neither can match the constraining power of multi-energy elastic data. Dispersive optical models [12] exploit precisely such energy correlations. The angle-resolved analysis (Fig. 7) shows

that backward-angle measurements ($\theta > 130^\circ$) are disproportionately valuable for breaking parameter degeneracies.

One caveat: the analysis uses the Woods-Saxon parameterization. However, the information limit arises from physics—elastic scattering probes an integrated “scattering surface” rather than the detailed radial shape—regardless of parameterization.

Conclusion.—This work provides the first quantitative decomposition of the constraining power of nucleon-nucleus scattering data on optical potential parameters. The step-by-step analysis reveals a clear hierarchy: (1) elastic angular distributions alone constrain only $D_{\text{eff}} \approx 1.7 \pm 0.4$ parameter combinations out of 11, primarily the potential radii; (2) the reaction cross section σ_R adds negligible information; (3) the analyzing power A_y lifts D_{eff} by $\sim 30\%$ through spin-orbit information; (4) multi-energy combination provides the largest improvement, reaching $D_{\text{eff}} \approx 2.2$; and (5) the 17-parameter KD02 global parametrization yields $D_{\text{eff}} \approx 1$ of 17—not because information is lost, but because KD02’s polynomial energy dependence (e.g., a cubic for $V(E)$) has more adjustable coefficients than the data can resolve.

Crucially, these limits are *scale-invariant*: improving experimental precision reduces error bars but cannot change the topology of the degenerate manifold. The result holds universally across the nuclear chart ($A = 12$ – 208 , $E = 10$ – 200 MeV, neutrons and protons). A global analysis combining all 168 systems in the 45-dimensional universal KD02 coefficient space yields $D_{\text{eff}} = 2.0/45$, with the dominant direction (69% of total information) being the real potential radius—confirming that the information limit is an intrinsic property of nucleon-nucleus scattering, not an artifact of analyzing a single system.

The “sloppy” parameter directions do not imply that these aspects of the potential are unconstrained by all data—rather, they are insensitive to single-energy scattering measurements. The multi-energy analysis (Fig. 4) demonstrates that combining data across energies probes orthogonal directions in parameter space, providing a constructive path forward. This supports the physical basis of dispersive optical models [12], which leverage the energy dependence of potentials as an additional constraint.

The angle-resolved analysis (Fig. 7) shows that backward-angle measurements ($\theta > 130^\circ$) are disproportionately valuable. For rare-isotope beam facilities where beam time is precious, the step-by-step decomposition provides actionable guidance: multi-energy angular distributions—even at moderate statistics—yield more constraining power than high-statistics measurements at a single energy, and A_y measurements are more informative than σ_R . The same framework could quantify parameter degeneracy in nucleon-nucleon potentials, suggesting that “sloppiness” may be a universal feature of effective nuclear interactions.

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* jinl@tongji.edu.cn

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End Matter

Relation to Bayesian analysis.—The Fisher information matrix (FIM) and Bayesian posterior are connected through a fundamental identity. For a likelihood $\mathcal{L}(\mathbf{d}|\boldsymbol{\theta})$ with flat priors, the posterior covariance in the Gaussian (Laplace) approximation is

$$\Sigma_{\text{post}} = F^{-1}, \quad F_{ij} = -\left\langle \frac{\partial^2 \log \mathcal{L}}{\partial \theta_i \partial \theta_j} \right\rangle. \quad (5)$$

The eigenvalues $\{\lambda_i\}$ of F are thus the inverse of the posterior variances along the principal axes: large λ_i corresponds to a tightly constrained direction (small posterior variance), while small λ_i corresponds to a sloppy direction (large posterior variance). The effective dimensionality $D_{\text{eff}} = (\sum \lambda_i)^2 / \sum \lambda_i^2$ counts the number of directions with comparable constraint.

This framework provides a rigorous explanation for the empirical findings of King *et al.* [6]. Their Bayesian MCMC analysis of nucleon- ^{48}Ca , ^{90}Zr , and ^{208}Pb elastic scattering (using a Woods-Saxon parameterization with 9 free central parameters, fixing spin-orbit) found that:

- Only the depth V and radius r of the real potential showed strong correlation in MCMC scatter plots;
- All other parameter pairs exhibited approximately circular (uncorrelated) scatter plots;
- The frequentist χ^2 approach produced spurious strong correlations among all parameters;

- The frequentist 95% confidence intervals were unrealistically narrow (44–92% empirical coverage versus the nominal 95%).

Our Fisher analysis explains each of these observations. The FIM eigenvector decomposition (Fig. 3) shows that \mathbf{e}_1 is dominated by r_v (65%) and V (12%), precisely the pair King *et al.* found correlated. The remaining parameters contribute to sloppy directions with eigenvalues 10^3 – 10^7 times smaller, rendering their correlations invisible in finite MCMC samples. The condition number exceeding 10^7 [Fig. 2(c)] explains why the frequentist approach—which approximates the posterior as a multivariate Gaussian around the χ^2 minimum—produced spurious correlations: projecting a highly anisotropic “cigar-shaped” posterior onto a Gaussian tangent plane creates artificial coupling between the elongated and compressed directions.

The narrow frequentist confidence intervals found in Ref. [6] (Table I therein) are also predicted by the Fisher analysis. The frequentist approach samples from the Gaussian approximation $\boldsymbol{\theta} \sim \mathcal{N}(\hat{\boldsymbol{\theta}}, F^{-1})$, which is valid only near the minimum. But with $D_{\text{eff}} \approx 1.7$, the χ^2 landscape has ~ 9 nearly flat directions where the Gaussian approximation breaks down, causing the frequentist method to dramatically underestimate the true parameter uncertainty along sloppy directions.

Computational advantages.—The Fisher approach offers several practical advantages over Bayesian MCMC for systematic studies:

1. *Efficiency*: Computing D_{eff} for one system requires $2N_p + 1 = 23$ evaluations of the scattering solver (for $N_p = 11$), taking ~ 20 seconds. King *et al.* required $\sim 100,000$ MCMC samples per system, with each sample requiring a FRESKO evaluation. The ~ 4000 -fold speedup enables the systematic scan across 168 configurations presented here—an analysis that would require $\sim 10^7$ total solver calls with MCMC.
2. *Analytical results*: The scale-invariance of D_{eff} (invariance under $F \rightarrow \alpha F$) is an exact mathematical property, whereas establishing this from MCMC would require running chains at multiple noise levels and comparing.
3. *Decomposability*: The Fisher matrix can be additively decomposed by observable ($F = F_{\text{elastic}} + F_{A_y} + F_{\sigma_R} + \dots$) and by energy ($F = \sum_E F(E)$), enabling the step-by-step analysis in Fig. 4. Bayesian posteriors cannot be similarly decomposed without re-running the full MCMC.
4. *Subgroup analysis*: Block sub-matrices of F quantify the information content for parameter subgroups (real volume, imaginary, spin-orbit), providing physical insight into which observables constrain which parameters.

The two approaches are complementary. The Fisher analysis efficiently maps the global structure of parameter constraints and proves fundamental properties (scale-invariance, universality), while Bayesian MCMC captures non-Gaussian features of the posterior—such as the asymmetric tails observed by King *et al.* [6]—that become important for precise credible intervals. A natural next step is to use the FIM eigenvectors to define an optimal reparameterization that separates stiff from sloppy directions, potentially improving MCMC convergence by orders of magnitude [9].

KD02 global parameter analysis.—The Koning-Delaroche potential [10] parametrizes the energy dependence of the 11 local parameters through 17 global parameters. For example, the real volume depth follows $V(E) = v_1[1 - v_2(E - E_f) + v_3(E - E_f)^2 - v_4(E - E_f)^3]$, where E_f is the Fermi energy. The geometry parameters (r_v , a_v , r_d , a_d) are energy-independent.

To analyze constraints on global parameters, I construct the Jacobian matrix $J = \partial \theta_{\text{local}} / \partial \theta_{\text{global}}$, where θ_{local} contains all $11 \times N_E$ local parameters across N_E energies. The global Fisher matrix is then $F_{\text{global}} = J^T F_{\text{combined}} J$, where $F_{\text{combined}} = \text{diag}(F(E_1), \dots, F(E_{N_E}))$.

With 7 energies and all observables, $D_{\text{eff}} \approx 1.0$ of 17 global parameters. The eigenvalue spectrum spans 15 orders of magnitude, with the largest eigenvalue corresponding to the energy-independent radius parameters.

As discussed in the main text, this low D_{eff} does not reflect a loss of constraining power relative to the multi-energy local analysis ($D_{\text{eff}} \approx 2.2$ of 11). The data still determine the same local-parameter combinations at each energy.

To understand why D_{eff} decreases when the parameter space grows, consider an analogy. Suppose a fixed “budget” of experimental information (say, 100 units) must be distributed among the parameters. In the 11-parameter local model, this budget is concentrated: ~ 2 parameters each receive ~ 45 units (well-constrained), while the remaining 9 share ~ 10 units (sloppy); the resulting $D_{\text{eff}} \approx 2$. Switching to the 17-parameter global model does not change the total budget, but introduces 6 additional parameters (the higher-order polynomial coefficients v_3 , v_4 , d_3 , etc.) that each receive effectively zero. These “dead” parameters dilute D_{eff} from $2.2/11$ to $1.0/17$ without affecting the actual constraints on any observable.

More concretely: $V(E) = v_1[1 - v_2(E - E_f) + v_3(E - E_f)^2 - v_4(E - E_f)^3]$ has four coefficients, but elastic data at 7 energies can only pin down v_1 (overall depth) and v_2 (linear energy slope). The difference between a quadratic and cubic energy dependence is far smaller than the uncertainty on V itself, so v_3 and v_4 are arbitrary. The effect is analogous to fitting a straight line with a fifth-order polynomial: the slope and intercept are well-determined, but the higher-order coefficients are unconstrained. This implies that future global optical potentials could use simpler energy-dependent functional forms without sacrificing predictive power.

Multi-system global analysis.—The per-nucleus analysis above uses 17 parameters specific to a given (A, Z) , since the KD02 formulas substitute nuclear properties into the coefficients. To test whether the information limit is truly universal, we extend the analysis to the full KD02 coefficient space: 45 universal parameters that generate all local potentials for any nucleus and energy. These include shared geometry coefficients (e.g., $r_v = r_{v0} - r_{v1}A^{-1/3}$ with $r_{v0} = 1.3039$, $r_{v1} = 0.4054$), shared depth amplitudes and energy scales, and separate neutron/proton depth coefficients.

For each of the 168 configurations, we compute the log-space Jacobian $J_i = \partial \log \theta_{\text{local}} / \partial \log \theta_{\text{universal}}$ and transform:

$$F_{\text{universal}} = \sum_{i=1}^{168} J_i^T F_i^{\text{local}} J_i. \quad (6)$$

The result is $D_{\text{eff}} = 2.0$ out of 45 universal parameters. The eigenvalue spectrum reveals a steep hierarchy: the first eigenvector (69% of total Fisher information) is dominated by the real potential radius coefficient r_{v0} ($|e_1| = 0.97$), confirming that the Igo ambiguity—the primacy of the potential radius as the one well-constrained quantity—persists as the defining feature even at the

global level. The second eigenvector (12%) corresponds to the surface imaginary radius r_{d0} . Together, the top 5 eigenvectors capture 97% of all information.

Crucially, D_{eff} saturates rapidly as nuclei are added: after combining ~ 5 (nucleus, projectile) systems, $D_{\text{eff}} \approx 2.2$; adding the remaining 19 systems provides no further improvement. This saturation demonstrates that different target nuclei probe essentially the same parameter combinations in the universal space—the constraining power of $n + {}^{40}\text{Ca}$ is not fundamentally different from that of $p + {}^{208}\text{Pb}$ when projected onto the universal coefficients.

Interestingly, elastic-only data yield $D_{\text{eff}} = 2.5/45$ in the global analysis, slightly *higher* than the all-observable result ($D_{\text{eff}} = 2.0/45$). This illustrates a key distinction: D_{eff} measures the *shape* of the eigenvalue spectrum (how uniformly information is distributed across directions), not the *total* information content ($\text{Tr } F =$

$\sum \lambda_i$, which always increases with more data). Adding A_y data increases the total Fisher information dramatically (individual eigenvalues grow by factors of 300–500), but the new information is deposited along already-stiff directions—predominantly the potential radius r_{v0} —rather than opening new sloppy directions. The eigenvalue spectrum becomes more peaked, and the participation ratio decreases. An analogy: illuminating an object with a brighter lamp from the same angle produces a sharper image of the same features, but does not reveal new sides; only changing the viewing angle—analogueous to varying the beam energy—can do that. This explains why multi-energy combination (Step 4) is the most effective route to increasing D_{eff} : different energies probe genuinely orthogonal parameter combinations. In both cases, only ~ 2 universal coefficient combinations are effectively constrained by all available nucleon-nucleus scattering data (Fig. 5).

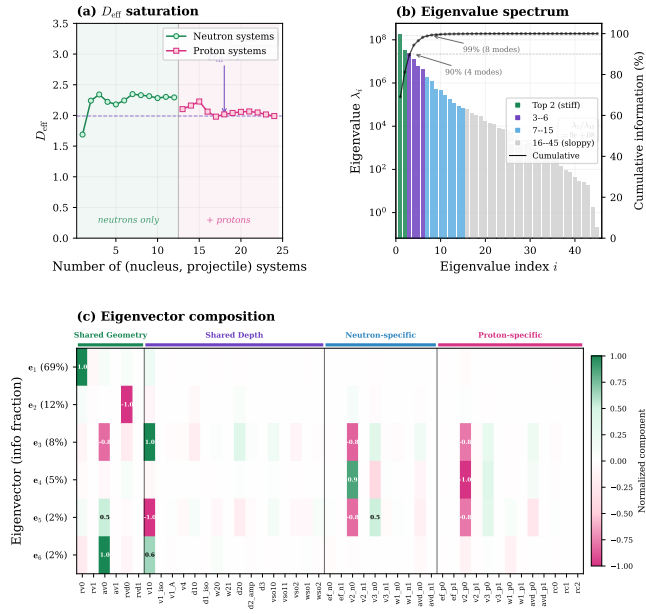


FIG. 5. Global Fisher analysis in the 45-dimensional universal KD02 coefficient space. (a) D_{eff} saturates at ~ 2 after combining ~ 5 systems; adding protons (pink) to neutrons (green) provides no improvement. D_{eff} can decrease when additional observables (e.g., A_y) primarily reinforce already-stiff directions, concentrating the eigenvalue spectrum (see text). (b) Eigenvalue spectrum showing the first mode carries 69% of all information; 4 modes suffice for 90%. (c) Eigenvector composition of the top 6 modes. The dominant mode e_1 is 97% r_{v0} (the real potential radius coefficient), confirming that the Igo ambiguity governs the global constraint structure.

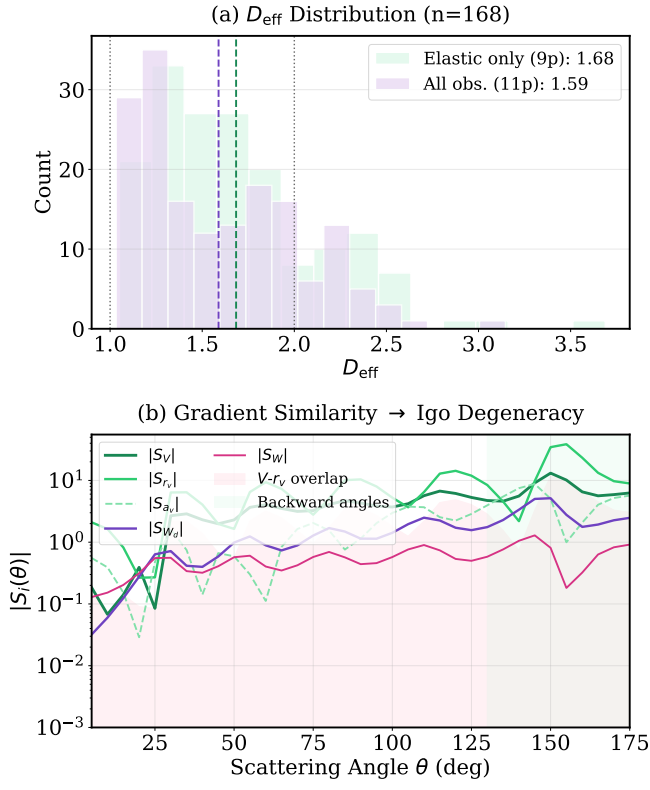


FIG. 6. Information geometry of the Igo ambiguity. The sensitivity vectors $|S_V(\theta)|$ and $|S_{r_v}(\theta)|$ have nearly identical angular dependence, pointing in the same direction in function space. This geometric near-collinearity is the origin of the depth-radius degeneracy.

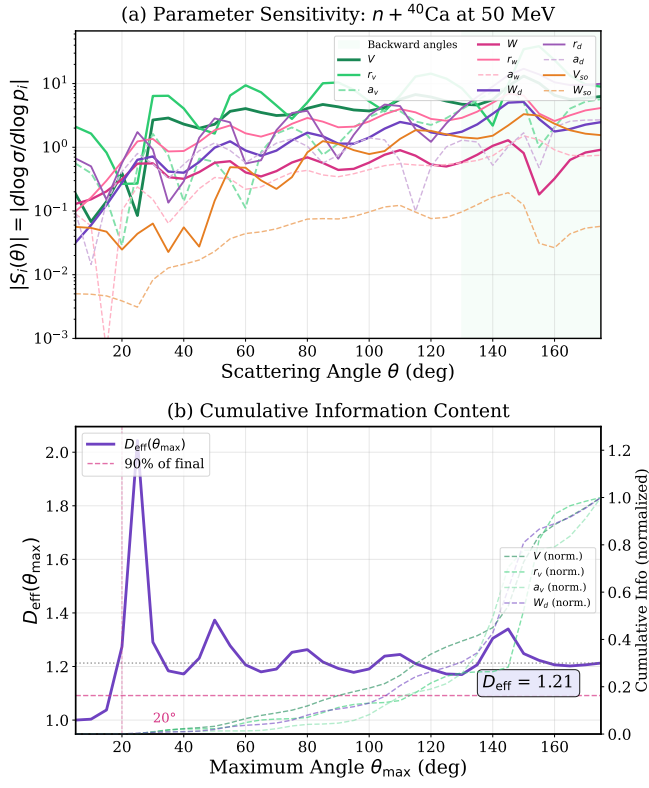


FIG. 7. Angle-resolved sensitivity for $n + {}^{40}\text{Ca}$ at 50 MeV. (a) Parameter sensitivity $|S_i(\theta)|$ showing near-identical angular patterns for V and r_v (Igo degeneracy); diffuseness parameters gain sensitivity at backward angles (shaded). (b) Cumulative $D_{\text{eff}}(\theta_{\text{max}})$ showing that $\sim 90\%$ of information content is captured by $\theta \approx 140^\circ$.