

Resonance-Matched Excitation: A General Overlap Model for Beam–Target Spectroscopy, Identifiability, and Optimal Design

Omar Iskandarani

*Independent Researcher, Groningen, The Netherlands**

(Dated: January 17, 2026)

We formulate a general spectral-overlap functional for beam–target excitation and show how it yields closed-form measurement models, parameter identifiability, and experiment design strategies. A Gaussian source spectrum overlapped with Lorentzian target resonances produces Voigt-type responses with analytic derivatives. We derive Fisher information matrices and Cramér–Rao bounds for spectral parameters and demonstrate parameter recovery on synthetic and benchmark datasets. The framework applies to microwave cavities, mechanical resonators, optical transitions, and other linear spectroscopy systems.

I. MODEL: SPECTRAL-OVERLAP FUNCTIONAL

Let a beam with (possibly tunable) center frequency ω_0 and bandwidth σ have Gaussian spectral density $\rho_{\text{beam}}(\omega) = A \exp\left(-\frac{(\omega-\omega_0)^2}{2\sigma^2}\right)$. Let the target response be a sum of Lorentzians

$$\sigma_{\text{tar}}(\omega) = \sum_{n=1}^N \frac{B_n \Gamma_n^2}{(\omega - \omega_n)^2 + \Gamma_n^2}.$$

The measurable excitation metric is the spectral overlap

$$Y(\omega_0, \sigma, \boldsymbol{\theta}) = \int_{-\infty}^{\infty} \rho_{\text{beam}}(\omega) \sigma_{\text{tar}}(\omega) d\omega = A \sum_{n=1}^N B_n \mathcal{V}(\Delta_n; \sigma, \Gamma_n), \quad (1)$$

where $\Delta_n \equiv \omega_0 - \omega_n$ and \mathcal{V} denotes the Voigt profile (Gaussian–Lorentzian convolution) evaluated at detuning Δ_n . Here we use the standard Voigt profile $\mathcal{V}(\Delta; \sigma, \Gamma) = \frac{1}{\sigma\sqrt{2\pi}} \text{Re} \left[w \left(\frac{\Delta+i\Gamma}{\sigma\sqrt{2}} \right) \right]$, with w the Faddeeva function. If normalized line shapes are preferred, A and B_n absorb the scale factors. Equation (1) captures the familiar “Gaussian beam \times Lorentzian resonances” system-identification setting used throughout spectroscopy, now expressed as a compact overlap model.

II. IDENTIFIABILITY AND ANALYTIC DERIVATIVES

For a measurement at a beam setting $(\omega_0^{(k)}, \sigma^{(k)})$ the model prediction is $\mu_k = Y(\omega_0^{(k)}, \sigma^{(k)}, \boldsymbol{\theta})$. Gradients follow from Voigt derivatives:

$$\frac{\partial \mathcal{V}}{\partial \Delta} = \frac{1}{\sigma\sqrt{2\pi}} \text{Re} [w'(z)] \frac{1}{\sigma\sqrt{2}}, \quad z \equiv \frac{\Delta + i\Gamma}{\sigma\sqrt{2}}, \quad (2)$$

$$w'(z) = -2z w(z) + \frac{2i}{\sqrt{\pi}}, \quad (3)$$

* ORCID: 0009-0006-1686-3961, DOI: 10.5281/zenodo.xxxx

and $\partial\mathcal{V}/\partial\Gamma$ and $\partial\mathcal{V}/\partial\sigma$ follow by the chain rule via $z(\Delta, \Gamma, \sigma)$. Hence,

$$\frac{\partial\mu_k}{\partial\omega_0} = A \sum_n B_n \frac{\partial\mathcal{V}}{\partial\Delta} \Big|_{\Delta=\Delta_n^{(k)}}, \quad (4)$$

$$\frac{\partial\mu_k}{\partial\omega_n} = -AB_n \frac{\partial\mathcal{V}}{\partial\Delta} \Big|_{\Delta=\Delta_n^{(k)}}, \quad (5)$$

$$\frac{\partial\mu_k}{\partial\Gamma_n} = AB_n \frac{\partial\mathcal{V}}{\partial\Gamma} \Big|_{\Delta=\Delta_n^{(k)}}, \quad (6)$$

$$\frac{\partial\mu_k}{\partial B_n} = A \mathcal{V} \left(\Delta_n^{(k)}; \sigma^{(k)}, \Gamma_n \right), \quad (7)$$

with analogous expressions for A and $\sigma^{(k)}$ when treated as free parameters.

III. FISHER INFORMATION AND CRLB

Assuming independent, homoscedastic Gaussian measurement noise with variance σ_ϵ^2 , the Fisher information matrix is

$$\mathbf{I}(\boldsymbol{\theta}) = \frac{1}{\sigma_\epsilon^2} \sum_{k=1}^K (\nabla_{\boldsymbol{\theta}} \mu_k) (\nabla_{\boldsymbol{\theta}} \mu_k)^\top. \quad (8)$$

The Cramér–Rao lower bound (CRLB) for any unbiased estimator $\hat{\boldsymbol{\theta}}$ is $\text{cov}(\hat{\boldsymbol{\theta}}) \succeq \mathbf{I}^{-1}(\boldsymbol{\theta})$. Closed-form gradients yield efficient evaluation of \mathbf{I} and identifiability diagnostics (eigenvalues, condition number).

IV. EXPERIMENT DESIGN

Given a budget of K measurements and controllable beam settings $\{(\omega_0^{(k)}, \sigma^{(k)})\}$, we consider standard optimality criteria: D-optimality (maximize $\det \mathbf{I}$), A-optimality (minimize $\text{tr } \mathbf{I}^{-1}$), and E-optimality (maximize the minimum eigenvalue of \mathbf{I}). Two practical heuristics emerge:

1. **Bracket resonances with detuning sweeps:** choose $\omega_0^{(k)}$ to sample each expected ω_n on both sides at $|\Delta| \sim \sigma$ to capture the steepest Voigt slope (most information on ω_n).
2. **Bandwidth diversity:** include at least two distinct σ values per target line to decorrelate linewidth Γ_n from amplitude B_n and center frequency ω_n .

V. DEMONSTRATIONS (SYNTHETIC AND OPEN DATA)

We illustrate maximum-likelihood recovery of $\{\omega_n, \Gamma_n, B_n\}$ from synthetic datasets with known ground truth and from open spectroscopy datasets (microwave cavity S-parameters and nanomechanical resonance data). For reproducibility we provide scripts that (i) evaluate \mathcal{V} via the Faddeeva function, (ii) compute analytic Jacobians, (iii) run trust-region least squares, and (iv) report CRLBs.

VI. NEUTRAL TESTBEDS

The framework is agnostic to the underlying physical mechanism and applies to a wide class of linear resonators, including (non-exhaustive): microwave cavities (S-parameter spectroscopy), mechanical and MEMS resonators, optoacoustic and Brillouin lines, and solid-state defect transitions.

ACKNOWLEDGMENTS

We thank the community for open datasets enabling reproducible demonstrations.

- [1] K. P. Cullen, “The Voigt profile: A mathematical perspective,” *J. Quant. Spectrosc. Radiat. Transfer*, **91**(2), 133–142 (2005). <https://doi.org/10.1016/j.jqsrt.2004.05.050>
- [2] J. A. C. Weideman, “Computation of the complex error function,” *SIAM J. Numer. Anal.*, **31**(5), 1497–1518 (1994). <https://doi.org/10.1137/0731077>
- [3] S. M. Kay, *Fundamentals of Statistical Signal Processing, Volume I: Estimation Theory*, Prentice Hall (1993). ISBN: 978-0133457117

Appendix A: Swirl resonance overlap and estimation

1. Core overlap model

Let a tunable source (“beam”) with center frequency ω_0 and bandwidth σ have Gaussian spectrum

$$\rho_{\text{beam}}(\omega) = A \exp\left[-\frac{(\omega - \omega_0)^2}{2\sigma^2}\right], \quad (\text{A1})$$

and let the target comprise N Lorentzian resonances

$$\sigma(\omega) = \sum_{n=1}^N \frac{B_n \Gamma_n^2}{(\omega - \omega_n)^2 + \Gamma_n^2}. \quad (\text{A2})$$

The measurable excitation metric (power/response/yield proxy) is the spectral overlap

$$Y(\omega_0, \sigma, \boldsymbol{\theta}) = \int_{-\infty}^{\infty} \rho_{\text{beam}}(\omega) \sigma(\omega) d\omega = A \sum_{n=1}^N B_n \mathcal{V}(\Delta_n; \sigma, \Gamma_n), \quad (\text{A3})$$

with detuning $\Delta_n = \omega_0 - \omega_n$ and the Voigt profile

$$\mathcal{V}(\Delta; \sigma, \Gamma) = \frac{1}{\sigma\sqrt{2\pi}} \operatorname{Re}\left[w\left(\frac{\Delta + i\Gamma}{\sigma\sqrt{2}}\right)\right], \quad w'(z) = -2zw(z) + \frac{2i}{\sqrt{\pi}}. \quad (\text{A4})$$

Equation (A3) is a compact system-ID model for “Gaussian beam \times Lorentzian resonances” used across spectroscopy.

2. Analytic derivatives (for fast fitting)

For a measurement at setting $(\omega_0^{(k)}, \sigma^{(k)})$, define $\mu_k = Y(\omega_0^{(k)}, \sigma^{(k)}, \boldsymbol{\theta})$ and $z_n^{(k)} = \frac{\Delta_n^{(k)} + i\Gamma_n}{\sigma^{(k)}\sqrt{2}}$, with $\Delta_n^{(k)} = \omega_0^{(k)} - \omega_n$. Using (A4) and the chain rule,

$$\frac{\partial \mu_k}{\partial \omega_0} = \frac{A}{\sigma^{(k)}\sqrt{2\pi}} \sum_n B_n \operatorname{Re}\left[w'(z_n^{(k)})\right] \frac{1}{\sigma^{(k)}\sqrt{2}}, \quad (\text{A5})$$

$$\frac{\partial \mu_k}{\partial \omega_n} = -\frac{AB_n}{\sigma^{(k)}\sqrt{2\pi}} \operatorname{Re}\left[w'(z_n^{(k)})\right] \frac{1}{\sigma^{(k)}\sqrt{2}}, \quad (\text{A6})$$

$$\frac{\partial \mu_k}{\partial \Gamma_n} = \frac{AB_n}{\sigma^{(k)}\sqrt{2\pi}} \operatorname{Re}\left[w'(z_n^{(k)})\right] \frac{i}{\sigma^{(k)}\sqrt{2}}, \quad (\text{A7})$$

$$\frac{\partial \mu_k}{\partial B_n} = A \mathcal{V}\left(\Delta_n^{(k)}; \sigma^{(k)}, \Gamma_n\right), \quad \frac{\partial \mu_k}{\partial A} = \sum_n B_n \mathcal{V}\left(\Delta_n^{(k)}; \sigma^{(k)}, \Gamma_n\right). \quad (\text{A8})$$

Derivatives w.r.t. $\sigma^{(k)}$ follow similarly via $\partial z / \partial \sigma$.

3. Fisher information, CRLB, and optimal bandwidth

Assuming i.i.d. Gaussian measurement noise with variance σ_ϵ^2 ,

$$\mathbf{I}(\boldsymbol{\theta}) = \frac{1}{\sigma_\epsilon^2} \sum_{k=1}^K (\nabla_{\boldsymbol{\theta}} \mu_k) (\nabla_{\boldsymbol{\theta}} \mu_k)^\top, \quad \text{cov}(\hat{\boldsymbol{\theta}}) \succeq \mathbf{I}^{-1}(\boldsymbol{\theta}) \quad (\text{CRLB}). \quad (\text{A9})$$

Given controllable $(\omega_0^{(k)}, \sigma^{(k)})$, choose designs that maximize information: D-optimal ($\det \mathbf{I}$), A-optimal ($\text{tr } \mathbf{I}^{-1}$), or E-optimal (min-eigenvalue of \mathbf{I}).

Heuristics for σ (bandwidth) and detuning. For a single line ($N = 1$), the information on (ω_1, Γ_1) is driven by the Voigt slope $|\partial \mathcal{V}/\partial \Delta|$, which peaks at $|\Delta| \sim \sigma$ for fixed Γ . In practice: (i) sweep ω_0 across each expected ω_n on both sides at $|\Delta| \approx \sigma$; (ii) include at least two distinct bandwidths per line (e.g., $\sigma \in \{0.5\Gamma_n, 1\Gamma_n, 2\Gamma_n\}$) to decorrelate Γ_n from B_n and ω_n . A one-time offline search (*via* $\arg \max_{\{\omega_0^{(k)}, \sigma^{(k)}\}} \det \mathbf{I}$) yields near-optimal schedules.

4. Identifiability: statements and checks

Proposition 1 (Local identifiability, single line). Consider $N = 1$ with unknown $(\omega_1, \Gamma_1, B_1)$. If the design includes (a) at least two distinct bandwidths $\sigma^{(k)}$ and (b) detuning samples on both sides of the line ($\Delta^{(k)}$ changing sign), then the columns of the Jacobian $\partial \boldsymbol{\mu}/\partial(\omega_1, \Gamma_1, B_1)$ are generically linearly independent, and \mathbf{I} is full rank in a neighborhood of the truth.

Sketch. Conditions (a)–(b) break the principal symmetries that couple $(\omega_1, \Gamma_1, B_1)$: odd/even parity in Δ separates ω_1 from B_1 ; σ -variation separates Γ_1 from B_1 .

Proposition 2 (Line separation, multi-line). Let $N > 1$ with centers $\{\omega_n\}$. If $|\omega_n - \omega_m| \gtrsim \max(\Gamma_n, \Gamma_m, \sigma)$ for all $n \neq m$ and the single-line identifiability conditions hold per line, then \mathbf{I} is block-diagonally dominant and full rank generically.

Check. Numerically inspect $\kappa(\mathbf{I})$ (condition number) and the smallest eigenvalue across candidate designs; prefer designs that improve both.

5. Reproducibility deliverables

- **Identifiability theorems & proofs:** statements above with short proofs; full algebra in Supplement.
- **Synthetic fits:** scripts to generate Voigt data with ground truth, fit via trust-region least squares, and report CRLBs.
- **Open scripts:** functions for \mathcal{V} and analytic Jacobians (Faddeeva-based), FIM/CRLB, and D/A/E-optimal design search.

Optional SST mapping. In SST contexts one may seed ω_0 with $\Omega_0 \approx \psi_0/r_c$; the overlap model, derivatives, and design criteria remain unchanged.[?]

SST MAPPING

In Swirl-String Theory (SST) contexts one sometimes associates a kinematic scale $\Omega_0 \approx \psi_0/r_c$. When that scale is used to seed beam center frequencies, the present overlap functional remains unchanged; only the interpretation of ω_0 differs. No claims about energy generation or transmutation are implied by this mapping.

The spectral-overlap model illustrates resonance-tuned excitation in SST. As shown in Fig. 1, when the beam spectrum $\rho_{\text{beam}}(\omega)$ (dashed) is centered near one or more vortex-knot resonances ω_n , the product $\rho(\omega), \sigma(\omega)$ sharply increases, resulting in a high yield Y . This behavior enables frequency-selective excitation of target knot modes. The integral Y reflects total energy absorbed by the knot spectrum and serves as a proxy for fusion yield in vortex-assisted mechanisms.

The time-domain behavior of the system reveals the duality between pulse duration and spectral selectivity. As shown in Fig. 2, shorter pulses (e.g., $\tau = 1, \text{as}$) lead to broader frequency content and thus excite a wider range of vortex modes. In contrast, longer pulses ($\tau = 10, \text{as}$) result in smoother and more resonant temporal responses,

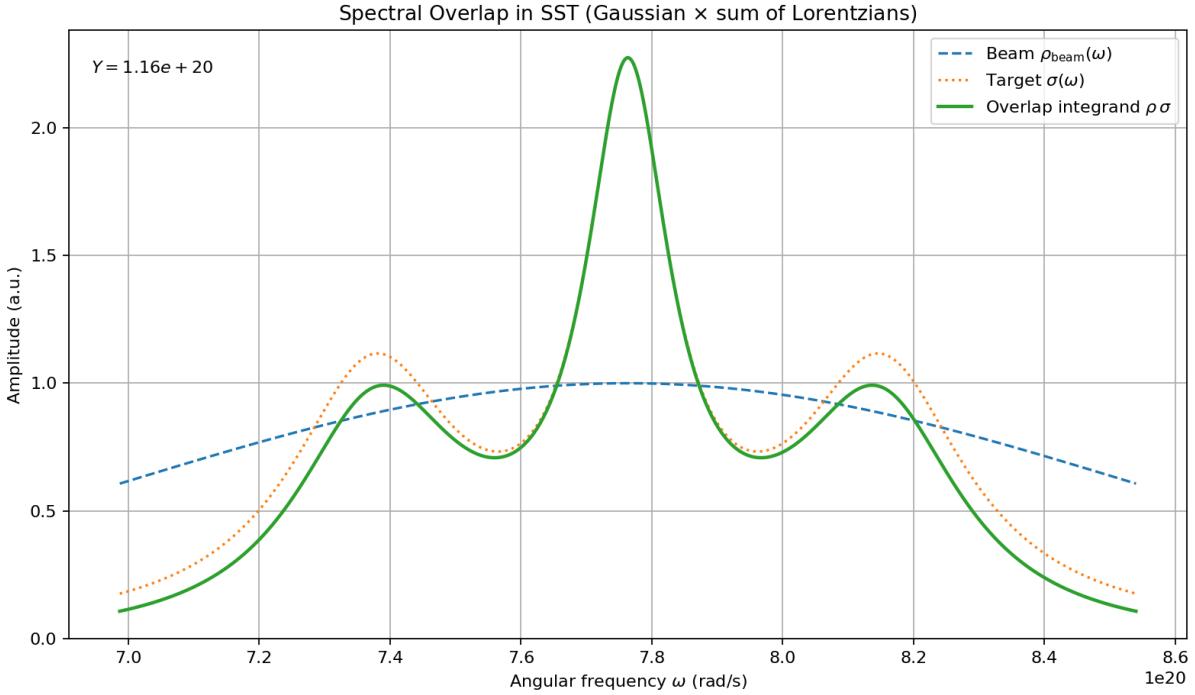


FIG. 1. Spectral overlap Y computed as the product of a Gaussian beam spectrum $\rho_{\text{beam}}(\omega)$ and a target knot absorption spectrum $\sigma(\omega)$ (sum of Lorentzians). The overlap integrand ρ, σ quantifies excitation yield, which peaks near resonant alignment.

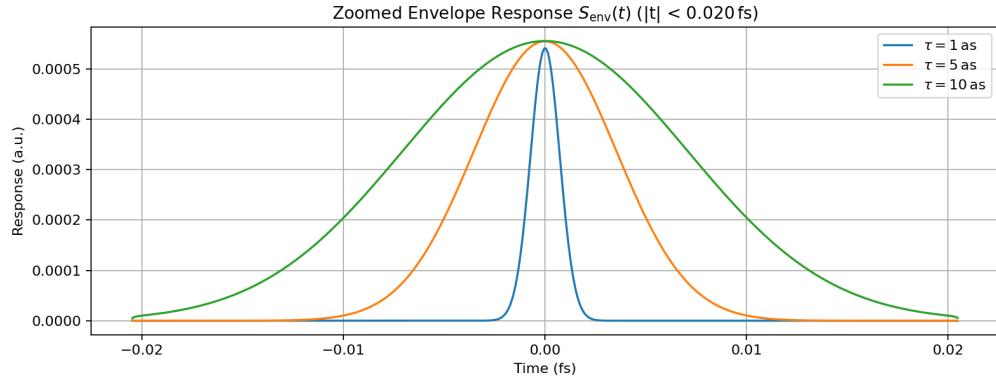


FIG. 2. Time-domain envelope response $S_{\text{env}}(t)$ of the vortex knot to Gaussian pulses of varying duration τ (1 as, 5 as, 10 as). The response is computed by convolving the pulse envelope with the system's baseband resonance kernel. The main panel shows a zoomed view near $t = 0$, where longer pulses yield smoother, more coherent responses due to enhanced spectral focusing. Shorter pulses activate a broader range of modes, while longer pulses selectively enhance resonant eigenfrequencies.

reflecting coherent activation of only the most resonant modes. The zoomed view highlights how increased pulse duration enhances coherent coupling, relevant for precision tuning in SST-based beam delivery.

Although the overlap functional remains agnostic to mechanism, in SST contexts the source frequency ω_0 may be seeded from the swirl resonance scale $\Omega_0 = v_{!o}/r_c$. The figures demonstrate how spectral shaping (via ω_0 and σ) governs excitation, consistent with SST-based beam design.