

Impact of resonance, Raman, and Thomson scattering on hydrogen line formation in Little Red Dots

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Accepted 2025 November 20. Received 2025 November 11; in original form 2025 August 11

ABSTRACT

Little Red Dots (LRDs) are compact sources at $z > 5$ discovered through *James Webb Space Telescope* spectroscopy. Their spectra exhibit broad Balmer emission lines ($\gtrsim 1000 \text{ km s}^{-1}$), alongside absorption features and a pronounced Balmer break – evidence for a dense, neutral hydrogen medium, in which the $n = 2$ state is significantly populated. When interpreted as arising from active galactic nucleus broad-line regions, inferred black hole masses from local scaling relations exceed expectations given their stellar masses, challenging models of early black hole–galaxy co-evolution. However, radiative transfer effects in dense media may also impact the formation of hydrogen emission lines. We model three scattering processes shaping hydrogen line profiles: resonance scattering by hydrogen in the $n = 2$ state, Raman scattering of ultraviolet (UV) radiation by ground-state hydrogen, and Thomson scattering by free electrons. Using 3D Monte Carlo radiative transfer simulations, we examine their imprint on line shapes and ratios. Resonance scattering produces strong deviations from Case B flux ratios, clear differences between $H\alpha$ and $H\beta$, and encodes gas kinematics in line profiles but cannot broaden $H\beta$ due to conversion to $\text{Pa}\alpha$. While Raman scattering can yield broad wings, scattering of the UV continuum is disfavoured given the absence of strong full width at half-maximum variations across transitions. Raman scattering of higher Lyman-series emission can produce $H\alpha/H\beta$ wing width ratios of $\gtrsim 1.28$, agreeing with observations. Thomson scattering can reproduce the observed $\gtrsim 1000 \text{ km s}^{-1}$ wings under plausible conditions – e.g. $T_e \sim 10^4 \text{ K}$ and $N_e \sim 10^{24} \text{ cm}^{-2}$ – and lead to black hole mass overestimates by factors $\gtrsim 10$. Our results provide a framework for interpreting hydrogen lines in LRDs and similar systems.

Key words: line: formation – radiative transfer – scattering – galaxies: active – galaxies: high-redshift – quasars: emission lines.

1 INTRODUCTION

Little Red Dots (LRDs) are compact sources in the early universe at $z > 5$ identified by *James Webb Space Telescope* (*JWST*) spectroscopy (Y. Harikane et al. 2023; D. D. Kocevski et al. 2023; H. B. Akins et al. 2025a; J. E. Greene et al. 2024; J. Matthee et al. 2024). They exhibit broad emission features with widths exceeding 1000 km s^{-1} in hydrogen and helium lines (M. Brazzini et al. 2025; I. Juodžbalis et al. 2025; V. Rusakov et al. 2025; A. Torralba et al. 2025; F. D’Eugenio et al. 2025c), typically interpreted as signatures of active galactic nuclei (AGNs). Many LRDs show a ‘v-shaped’ spectrum with a blue ultraviolet (UV) continuum and a red UV to optical slope (e.g. H. B. Akins et al. 2025a; X. Ji et al. 2025a), often a Balmer break (D. J. Setton et al. 2025; B. Wang et al. 2024). While LRDs often lack traditional AGN indicators, such as strong X-ray (A. Sacchi & A. Bogdan 2025; R. Maiolino et al. 2025b), radio emission (A. J. Gloude-mans et al. 2025), or an infrared (IR) excess (C. C. Williams et al. 2023; M. Xiao et al. 2025), other faint

indicators such as Fe II emission (E. Lambrides et al. 2024) or high ionization emission lines (M. Tang et al. 2025, cf. E. Lambrides et al. 2024) have been identified. This rather unusual combination of spectral properties challenges standard AGN interpretations and the applicability of standard calibrations of bolometric luminosity.

LRDs appear to host unexpectedly massive supermassive black holes (SMBHs, F. Pacucci et al. 2023; K. Inayoshi 2025). The broad emission features in $H\alpha$ and $H\beta$ allow estimates of black hole mass M_{BH} using single epoch virial methods (J. E. Greene & L. C. Ho 2005; A. E. Reines & M. Volonteri 2015; J.-H. Woo et al. 2015), which are commonly applied to nearby AGNs. Using this approach, recent studies (I. Juodžbalis et al. 2024; K. Inayoshi & R. Maiolino 2025; I. Juodžbalis et al. 2025; A. J. Taylor et al. 2025a, b) have estimated $M_{\text{BH}} \sim 10^{8-9} M_{\odot}$ for LRDs. However, it remains unclear whether Balmer line widths are reliable tracers of M_{BH} in these systems (see also lower M_{BH} by X-ray; T. T. Ananna et al. 2024). V. Rusakov et al. (2025) suggest that the broad wings are shaped by electron (Thomson) scattering, based on their observed exponential profiles, and derive SMBH masses that are $\sim 10 - 100$ times smaller when using Balmer line luminosities instead of widths. Additionally, R. P. Naidu et al. (2025) propose that resonance scattering of $H\beta$

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photons in a dense neutral medium contributes to the line broadening, offering an alternative explanation to a virialized broad-line region (BLR). Thus, while potentially crucial radiative transfer (RT) effects in LRDs are discussed in the literature, a systematic theoretical study investigating those effects is still missing.

Spectra of LRDs frequently exhibit not only broad Balmer emission lines but also complex profiles, including absorption features, P-Cygni-like shapes in the $H\alpha$ and $H\beta$ lines, and significant Balmer breaks. These features point to the presence of gas that is optically thick to Balmer transitions, likely due to an abundant population of hydrogen atoms in the $n = 2$ state (I. Juodžbalis et al. 2024; J. Matthee et al. 2024; A. Graaff et al. 2025; K. Inayoshi & R. Maiolino 2025; R. P. Naidu et al. 2025; F. D'Eugenio et al. 2025a, b; X. Ji et al. 2025a; R. Maiolino et al. 2025a). Interestingly, $H\alpha$ and $H\beta$ often show distinct line profiles: $H\alpha$ frequently shows a P-Cygni-like shape with blueshifted absorption and a redshifted emission peak, while $H\beta$ tends to exhibit deeper absorption and a less prominent red peak. In contrast, Paschen lines typically lack absorption features (M. Brazzini et al. 2025; I. Juodžbalis et al. 2025). These observational trends suggest that RT effects (i.e. resonance scattering of Balmer lines) may play a key role in shaping the emergent line profiles.

While resonance scattering is typically invoked to explain the frequency and spatial diffusion of Lyman- α ($Ly\alpha$) in neutral hydrogen environments (D. A. Neufeld 1990; M. Dijkstra et al. 2019), similar processes can occur for Balmer lines in environments like LRDs, where a significant $n = 2$ population exists. Thus, modeling the resonance scattering of Balmer photons is crucial for understanding hydrogen line formation in these systems. However, importantly, the atomic physics of Balmer resonance scattering differs from that of $Ly\alpha$. For example, electrons in the atomic hydrogen excited by $H\beta$ photons from $n = 2$ to 4 can decay via multiple de-excitation channels to $n = 3$, 2, or 1, with each branch leading to different line formations, while $H\alpha$ involves transitions from $n = 3$ to 2. As a result, $H\beta$ photons are more likely to be converted into $P\alpha$ via cascades to the $n = 4 \rightarrow 3$ transition, due to multiple scatterings. These distinctions highlight the necessity for RT models that incorporate both these multiple branches of Balmer line scattering and the physical and kinematic properties of the scattering medium.

Other scattering processes can also produce broad hydrogen line profiles. Thomson scattering by free electrons can generate symmetric, exponential wings around emission lines, with widths scaling with the electron thermal speed ($\sim 500 \text{ km s}^{-1}$ at $T = 10^4 \text{ K}$). Such wings have been observed and studied in Type IIIn supernovae (L. Dessart et al. 2009), symbiotic stars (M. Sekeráš & A. Skopal 2012; S.-J. Chang et al. 2018b), AGNs (R. J. Weymann 1970; R. A. Sunyaev & L. G. Titarchuk 1980; H.-W. Lee 1999; A. Laor 2006), and recently in LRDs (V. Rusakov et al. 2025 cf. M. Brazzini et al. 2025).

Raman scattering, where UV photons near Lyman-series wavelengths are inelastically scattered by ground-state hydrogen atoms, can also generate extremely broad, often asymmetric features around Balmer lines like $H\alpha$ and $H\beta$ (S.-J. Chang et al. 2015). Raman wings have been observed and investigated in symbiotic stars (H. Nussbaumer, H. M. Schmid & M. Vogel 1989; J. J. Yoo, J.-Y. Bak & H.-W. Lee 2002; S.-J. Lee & S. Hyung 2018; S.-J. Chang et al. 2018b), dense star-forming regions (M. A. Dopita et al. 2016; W. J. Henney 2021), and AGNs (S.-J. Chang et al. 2015; M. Kokubo 2024), and are proposed for LRDs as well (M. Kokubo & Y. Harikane 2025). These scattering mechanisms offer alternative explanations for the broad Balmer components in LRDs, independent of virialized motion in a classical AGN BLR. Even if scattering is not the dominant mechanism, its contribution complicates the interpretation of line widths and, therefore, affects standard black hole mass estimation

methods, as the observed broadening may additionally originate from scattering processes rather than purely kinematic broadening.

Given the complex line shapes and multiple broadening mechanisms, we investigate how RT effects influence hydrogen emission-line profiles in LRDs, focusing on $H\alpha$ and $H\beta$. We perform Monte Carlo RT simulations including three key processes: resonance scattering of $H\alpha$ and $H\beta$, Thomson scattering by free electrons, and Raman scattering by neutral hydrogen. Section 2 details the physics of each mechanism, the model geometry, and the simulation setup. Section 3 presents the emergent spectra and explores how each process shapes the line profiles. In Section 4, we examine observable signatures such as line profiles and explore the impact of scattering on the estimation of black hole mass.

2 METHOD

To understand the formation of Balmer line profiles in the LRDs, we conduct Monte Carlo RT simulations focusing on three scattering processes: (i) resonance scattering of Balmer lines, $H\alpha$ and $H\beta$, by hydrogen atoms in the metastable $2s$ state, (ii) Raman scattering of UV continuum photons by ground-state hydrogen atoms, and (iii) Thomson scattering of Balmer line photons by free electrons. Each mechanism can contribute to the broadening and shaping of emission lines, producing features such as P-Cygni profiles, double peaks, or broad wings in the spectra.

In order to explore the impact of scattering processes on observables, our simulations adopt a simplified spherical geometry, where a central point source is embedded within a uniform scattering medium. For each process, we use appropriate Monte Carlo RT codes to track photon propagation and scattering interactions, accounting for the specific cross-sections, branching ratios, and physical conditions relevant to each scattering mechanism. The following sections describe the setup, physics, and simulation methods for each scattering process in detail.

2.1 Radiative transfer framework and geometry

To investigate the formation of hydrogen emission lines in LRDs, we consider RT through different scattering processes within a simplified geometry. Our model consists of a central isotropic point source embedded in a uniform scattering medium, as illustrated in Fig. 1. The central source emits either hydrogen line photons (for resonance and Thomson scattering) or a flat UV continuum (for Raman scattering).

The geometry of the scattering medium is spherical, characterized by an outer radius R and an inner radius set at $0.01R$. The number density of hydrogen atoms or electrons is uniform within this region. For each scattering mechanism, we apply relevant physical parameters that control the resulting line profiles:

(i) For *resonance scattering* of Balmer lines, $H\alpha$ and $H\beta$, we consider the neutral hydrogen column density in the $2s$ state, $N_{\text{HI},2s}$, the random velocity dispersion σ_R , and an optional bulk radial outflow velocity v_{exp} . The velocity field parameters σ_R and v_{exp} influence the frequency shifts and the formation of double-peaked profiles or P-Cygni features.

(ii) For *Raman scattering* of the UV continuum, we adopt a flat incident spectrum covering the Lyman-series region from $Ly\beta$ to $Ly\delta$. The process depends primarily on the total H I column density N_{HI} , as the broadening of Raman-scattered features is dominated by the interaction cross-section, branching probabilities near the Lyman transitions, and the energy conservation during Raman scattering

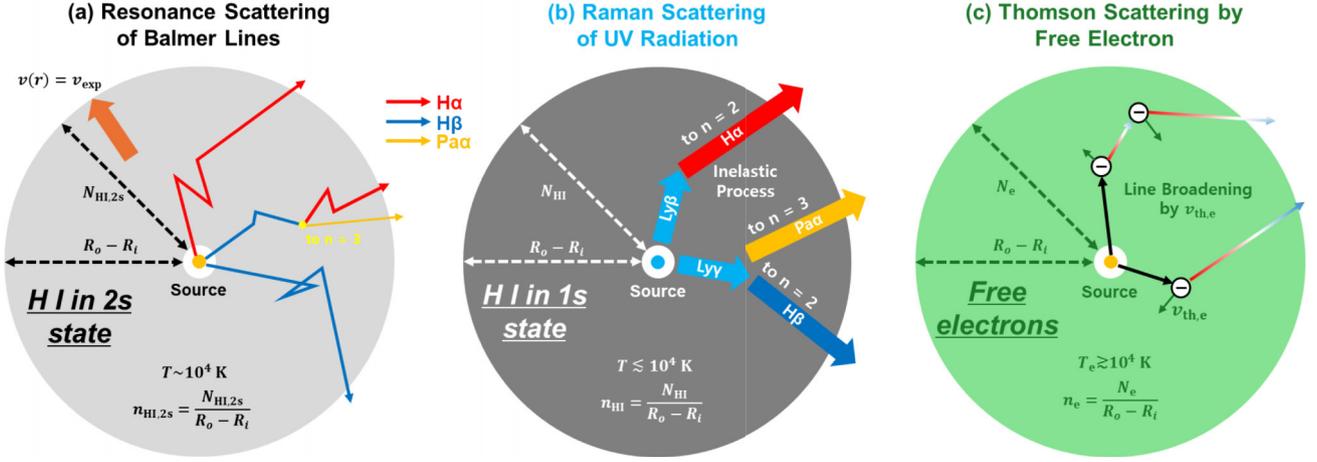


Figure 1. Schematic illustration of the model geometry. A central point source emits either $H\alpha$ and $H\beta$ photons (orange points for resonance and Thomson scattering) or a UV continuum (a blue point for Raman scattering), surrounded by a spherical scattering region. The sphere is divided into three wedges representing the different scattering processes: (a) resonance scattering of Balmer lines, $H\alpha$ and $H\beta$, in a spherical H I region populated with $2s$ -state hydrogen atoms, characterized by a column density $N_{\text{HI},2s}$, a random motion σ_{R} , and an expansion velocity v_{exp} . (b) Raman scattering of UV photons near the Lyman series by ground-state H atoms with a column density N_{HI} , producing broad optical and IR features (e.g. $H\alpha$, $H\beta$, and $\text{Pa}\alpha$). (c) Thomson scattering of Balmer line photons by free electrons in a spherical H II region, characterized by electron temperature T_e and electron density n_e , producing symmetric broadening proportional to the electron’s thermal speed $v_{\text{th},e}$.

Table 1. Key parameters for each scattering process. The ranges represent typical values used in our simulations.

Type of scattering	Parameter	Typical range
Resonance	H I column density in the $2s$ state $N_{\text{HI},2s}$	$10^{12} - 10^{16} \text{ cm}^{-2}$
	Expansion velocity v_{exp}	$0 - 500 \text{ km s}^{-1}$
	Random motion of H I gas σ_{R}	13.6 (hydrogen thermal motion at $T = 10^4 \text{ K}$) $- 100 \text{ km s}^{-1}$
	Width of intrinsic emission σ_{src}	$0 - 400 \text{ km s}^{-1}$
Raman	H I column density N_{HI}	$10^{20} - 10^{23} \text{ cm}^{-2}$
Thomson	Electron temperature T_e	$10^4 - 10^6 \text{ K}$
	Thomson optical depth τ_{Th}	$0.01 - 10$
	Width of intrinsic emission σ_{src}	$50 - 500 \text{ km s}^{-1}$

(i.e. Raman broadening). We neglect the effects of bulk or turbulent motions, as they have a negligible impact on the broadening of Raman features compared to the intrinsic energy shift from the inelastic scattering.

(iii) For *Thomson scattering* of Balmer lines, the scattering region is an ionized H II sphere characterized by the electron density n_e , temperature T_e , and the resulting Thomson optical depth $\tau_{\text{Th}} = \sigma_{\text{Th}} n_e R$. The broadening of the scattered line profiles arises primarily from the thermal motions of electrons $v_{\text{th}} = \sqrt{2kT/m_e} \approx 548 \text{ km s}^{-1}$ at $T_e = 10^4 \text{ K}$. Similar to Raman scattering, we neglect the effect of bulk outflows or turbulence, as these do not significantly impact the Thomson-scattering process due to the high thermal speed of electrons and wavelength-independent scattering cross-section.

The RT simulations for each process are performed using dedicated Monte Carlo codes, as detailed in Sections 2.2–2.4. The input parameters for each scattering process are summarized in Table 1.

2.2 Resonance scattering of Balmer lines

Resonance scattering in Balmer lines is expected in environments where the $n = 2$ population of hydrogen is significantly enhanced. In LRDs, the presence of Balmer absorption features and a strong Balmer break is indicative of this condition. Since $H\alpha$ and $H\beta$

photons can be trapped and experience multiple scatterings within optically thick gas, their emergent profiles are sensitive to the velocity structure and geometry of the medium. Modelling this process is essential to explain their P-Cygni-like profiles and the asymmetric absorption observed in Balmer lines, particularly the different behaviours of $H\alpha$ and $H\beta$.

2.2.1 Physical mechanism

Resonance scattering of Balmer lines occurs when hydrogen atoms in the $2s$ state absorb photons near the $H\alpha$ and $H\beta$ transition energies, exciting electrons to the $n = 3$ or 4 states, respectively. The subsequent de-excitation of the electron can result in the re-emission of a photon at the same wavelength or, through branching, produce photons in other lines such as $\text{Pa}\alpha$. This process is analogous to $\text{Ly}\alpha$ scattering (see e.g. review by M. Dijkstra et al. 2019), but for Balmer transitions.

The key role of the $2s$ state arises from its long lifetime (approximately 0.1 s), which is orders of magnitude longer than the lifetimes of the np excited states ($\sim 10^{-9}$ s). This long lifetime enables a significant population of hydrogen atoms to accumulate in the $2s$ state, providing a reservoir for resonance scattering. The atomic structure of hydrogen and relevant transitions are illustrated in Fig. 2.

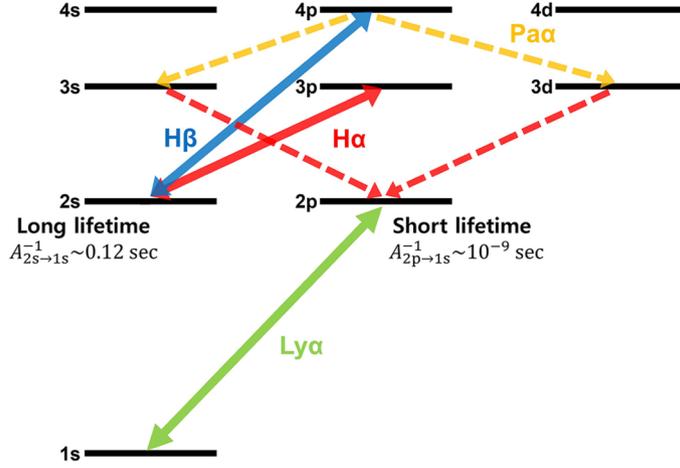


Figure 2. Energy levels of atomic hydrogen in $n = 1 - 4$. Solid arrows indicate the ‘resonant transitions’ involved in resonance scattering (see Section 2.2), while dashed arrows show other possible transitions for $H\alpha$, $H\beta$, and $Pa\alpha$ (omitting the higher Lyman transitions; ‘Case B’ assumption).

Table 2. Atomic parameters of $H\alpha$ and $H\beta$ to calculate their scattering cross-section in equation (1).

	$H\alpha(2s \rightarrow 3p)$	$H\beta(2s \rightarrow 4p)$
f	0.4360	0.1028
Γ	$4.4 \times 10^7 \text{ s}^{-1}$	$8.4 \times 10^6 \text{ s}^{-1}$
ν_0	$4.5600 \times 10^{14} \text{ s}^{-1}$	$6.1694 \times 10^{14} \text{ s}^{-1}$

Notes. The first row shows the oscillator strength from $2s$ to np state. The second and third rows represent the damping term and line centre frequency.

The scattering cross-section for a Balmer line is described by a Voigt profile:

$$\sigma_v = \frac{\sqrt{\pi} e^2}{m_e c} \frac{f}{\Delta \nu_D} H(a, x), \quad (1)$$

where f is the oscillator strength, $\Delta \nu_D$ is the Doppler width, and $H(a, x)$ is the Voigt function:

$$H(a, x) = \frac{a}{\pi} \int_{-\infty}^{\infty} \frac{e^{-y^2}}{(x-y)^2 + a^2} dy. \quad (2)$$

Here, $a = \Gamma/(4\pi\Delta\nu_D)$ is the natural width parameter, and $x = (\nu - \nu_0)/\Delta\nu_D$ is the dimensionless frequency offset. The thermal Doppler width is $\Delta\nu_D = \nu_0 v_{th}/c$, where $v_{th} = \sqrt{2kT/m_H}$ is the thermal speed of hydrogen at a temperature T . As the total random motion $\sigma_R = \sqrt{v_{th}^2 + v_{turb}^2}$, assuming the microturbulent approximation with the turbulent speed v_{turb} , we set the thermal speed v_{th} to be the random motion of hydrogen σ_R in our simulation.

Adopting the atomic data in Table 2, the optical depths at line centre for $H\alpha$ and $H\beta$ are given by:

$$\tau_{0,H\alpha} \approx 33.4 \left(\frac{N_{HI,2s}}{10^{14} \text{ cm}^{-2}} \right) \left(\frac{\sigma_R}{12.8 \text{ km s}^{-1}} \right)^{-1}, \quad (3)$$

$$\tau_{0,H\beta} \approx 5.83 \left(\frac{N_{HI,2s}}{10^{14} \text{ cm}^{-2}} \right) \left(\frac{\sigma_R}{12.8 \text{ km s}^{-1}} \right)^{-1}. \quad (4)$$

2.2.2 Branching ratios and line conversion

When an $H\alpha$ photon excites a $2s$ electron to the $3p$ state, the electron can decay to either the $n = 1$ ($1s$) or $n = 2$ ($2s$) state, emitting $Ly\beta$ or $H\alpha$, respectively (see also Fig. 2). The branching ratio to the $1s$ state is significantly larger (0.882) than that to the $2s$ state

Table 3. Branching ratio of permitted transitions at $3p$ and $4p$ states.

Initial state	$3p$	$3p$	$4p$	$4p$	$4p$
Final state	$1s$	$2s$	$1s$	$2s$	$3s, 3d$
Branching ratio	0.882	0.118	0.839	0.119	0.042
Branching ratio w/o $1s$	–	1	–	0.739	0.261

(0.118), as shown in Table 3. However, under the Case B assumption, the $3p \rightarrow 1s$ transition is neglected, as $Ly\beta$ photons are typically reabsorbed by the surrounding HI medium with the $1s$ state. As the population of the $n = 1$ state is much higher than that of the $n = 2$ state, if the gas is optically thick for Balmer lines, it is also optically thick for Lyman series. Thus, the Case B assumption is natural in the scattering medium for resonance scattering. Consequently, $H\alpha$ photons predominantly remain $H\alpha$ throughout the scattering process.

For $H\beta$ photons, excitation to the $4p$ state allows decay to the $n = 1$ ($1s$), $n = 2$ ($2s$), or $n = 3$ ($3s/3d$) states. The raw branching ratios for these transitions are 0.839, 0.119, and 0.042, respectively. When the $1s$ transition is suppressed under the Case B assumption, the relative branching ratio to the $n = 3$ state increases to 0.261. Therefore, the conversion of $H\beta$ photons into $Pa\alpha$ via the $4p \rightarrow 3s/3d$ transition is not negligible and must be considered in the simulation for the optically thick medium. This process also generates $H\alpha$ photons at the scattering location, as the $Pa\alpha$ cascade involves a subsequent $3s/3d \rightarrow 2p$ transition (see Fig. 2).

2.2.3 Monte Carlo simulation

The resonance scattering simulations are performed using the 3D Monte Carlo code RT-SCAT (S.-J. Chang et al. 2023a; S.-J. Chang & M. Gronke 2024) in which discrete photon packages are followed through real and frequency space until they escape the simulation domain, at which point the emergent properties are recorded. Specifically, we follow these steps:

- (i) Define the spherical geometry characterized by an HI column density in the $2s$ state $N_{HI,2s}$, a random motion σ_R , and an expansion velocity v_{exp} ; the radial bulk motion $v(r)$ is constant (i.e. $v(r) = v_{exp}$).
- (ii) Generate isotropic $H\alpha$ and $H\beta$ photons from the central source, with an intrinsic Gaussian width σ_{src} .

(iii) For each photon, determine the scattering location. Using a dimensionless free path of the photon, $\tau_f = -\ln r$, where r is a random number drawn uniformly in $[0,1]$, a free path of a photon l_f is given by τ_f/σ_v ; σ_v is a scattering cross-section (equation 1). The frequency ν is adopted as the frequency in the rest frame of the cell, where the photon penetrates.

(iv) Determine the scattering outcome:

(a) For $H\alpha$, photons always remain $H\alpha$ after scattering.

(b) For $H\beta$, the final state is selected by the branching ratios. If a random number r drawn uniformly in $[0,1]$ is smaller than the branching ratio to the $n = 2$ state, 0.739, the photon remains $H\beta$. Otherwise, it converts to $Pa\alpha$ with an accompanying $H\alpha$ photon at the same location (with the wavelength at the line centre in the atom rest frame). Because of the low number of H I atoms in the $n = 3$ state, the $Pa\alpha$ photon immediately escapes, while the newly generated $H\alpha$ photon continues to propagate in the spherical H I region.

(v) Compute the new photon direction with the isotropic distribution and frequency shift by a random motion σ_R . While the incident and scattered frequencies are identical in the rest frame of atomic hydrogen, which scatters a photon, the scattered frequency in the rest frame of the cell is changed due to the Doppler shift caused by the random motion of the hydrogen.

(vi) Repeat steps (iii)–(v) until the photon escapes the grid. When collecting photons, the photon's frequency is converted to the frequency in the observed frame, taking into account the kinematics at the last scattering location.

(vii) Back to a step (ii) until the total number of intrinsic photons becomes 10^8 .

A visual representation of this process is shown in the left panel of Fig. 1. The detailed Monte Carlo method of RT-SCAT is described in section 2.5 of S.-J. Chang & M. Gronke (2024). We will show simulated results of resonance scattered $H\alpha$, $H\beta$, and $Pa\alpha$ profiles in Section 3.1.

2.3 Raman scattering

Raman scattering generates extremely broad wings around Balmer lines, due to the inelastic scattering of UV photons near Lyman-series wavelengths by neutral hydrogen. While this process requires strong UV radiation fields and large N_{HI} , both conditions are plausible in LRDs, which show V-shaped spectral energy distributions (SEDs) and deep Balmer breaks. If Lyman emission lines or a strong UV continuum exist, Raman scattering can significantly reshape the Balmer line profiles and even mimic AGN-like broad components.

2.3.1 Physical mechanism

Raman scattering with H I atom is an inelastic process in which a UV photon near the hydrogen Lyman series is scattered by a ground-state hydrogen atom, producing an optical or IR photon through de-excitation to $n > 1$ states. If the electron returns to the ground state ($n = 1$) after scattering, the process is elastic, and the photon retains its original energy; this is known as Rayleigh scattering. In contrast, if the electron de-excites to higher energy levels ($n > 1$), the scattered photon undergoes an energy shift corresponding to the difference between the Lyman transition energy and the new state. Specifically, UV continuum photons near $Ly\beta$ form scattered features around $H\alpha$ through Raman scattering to the $n = 2$ state. Fig. 3 illustrates how Raman scattering of an $Ly\gamma$ photon to $n = 2$ and 3 states results in the conversion of the UV photon to $H\alpha$ and $Pa\alpha$, respectively.

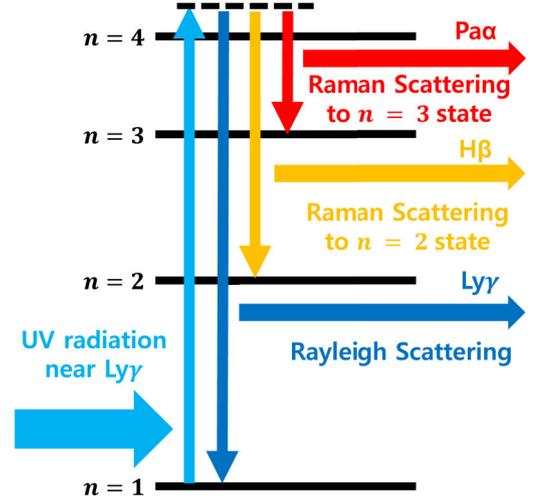


Figure 3. Schematic illustration of Raman scattering of UV radiation near $Ly\gamma$. The black solid lines represent the energy levels of atomic hydrogen. The light blue line indicates the incident UV radiation near $Ly\gamma$. The blue arrow represents Rayleigh scattering, in which the electron de-excites to the $n = 1$ state, emitting $Ly\gamma$. The red and orange arrows represent Raman scattering to the $n = 2$ and 3 states, producing $H\beta$ and $Pa\alpha$, respectively.

Computing the probability of Raman scattering occurring as a function of a wavelength in the far-UV regime is non-trivial. While the scattering probability for Rayleigh and Raman scatterings is given by the total cross-section, the individual transition probabilities are given by the branching ratios, which are the ratio between the cross-section for Raman scattering divided by the total scattering cross-section. Thus, the sum of the branching ratios to the $n = 1$ state (Rayleigh scattering) and $n > 1$ states (Raman scattering) is unity. Fig. A1 shows the total scattering cross-section and branching ratios to each n state.

This process involves an energy shift corresponding to the Lyman transition energy, given by

$$E_{\text{inc}} = E_{\text{sc}} + E_{\text{Lyman}}, \quad (5)$$

where E_{inc} is the energy of the incident UV photon, E_{sc} is the energy of the scattered photon, and E_{Lyman} is the energy of the corresponding Lyman transition. For instance, following the example to $Pa\alpha$ depicted in Fig. 3 (light blue to red arrow), $E_{\text{inc}} \approx E_{\text{Ly}\gamma}$, $E_{\text{Lyman}} = E_{\text{Ly}\beta}$, and $E_{\text{sc}} \approx E_{\text{Pa}\alpha}$.

The scattered wavelength is computed correspondingly as

$$\frac{1}{\lambda_{\text{sc}}} = \frac{1}{\lambda_{\text{inc}}} - \frac{1}{\lambda_{\text{Lyman}}}, \quad (6)$$

where λ_{Lyman} is the wavelength of the relevant Lyman transition (e.g. $Ly\beta$ and $Ly\gamma$).

This scattering process leads to an emergent broad line. To first order, the emergent line shape follows the corresponding cross-section (cf. Fig. A1). However, the broadening of Raman-scattered features arises not only from the natural cross-section dependence but also from the energy shift. This can be illustrated by the differential relation between the incident and scattered wavelengths, which is given by

$$d\lambda_{\text{sc}} = \left(\frac{\lambda_{\text{sc}}}{\lambda_{\text{inc}}} \right)^2 d\lambda_{\text{inc}}, \quad (7)$$

where the factor $(\lambda_{\text{inc}}/\lambda_{\text{sc}})^2$ represents the additional broadening from the Raman-scattering process. As the derivative of the velocity is given by $dv = d\lambda/\lambda$, this factor becomes $\lambda_{\text{sc}}/\lambda_{\text{inc}}$ in velocity space, which is ~ 6.4 and 5 for $H\alpha$ and $H\beta$, respectively.

Due to the cross-section dependence, the overall broadening of Raman features depends primarily on the total HI column density N_{HI} , with higher N_{HI} leading to broader features with widths exceeding several thousand km s^{-1} (H.-W. Lee & J.-H. Yun 1998; S.-J. Chang et al. 2015, 2018b; M. Kokubo 2024). Unlike resonance scattering, the effects of the bulk velocity v_{exp} or the random motion σ_{R} are negligible compared to the intrinsic broadening from the energy shift, assuming the continuum radiation as an input source. However, if UV emission lines, such as He II lines at 1025, 972, and 949 Å and O VI resonance lines at 1032 and 1038 Å (rather than the continuum) undergo Raman scattering, the kinematics of the HI region and the profile of input radiation (H. M. Schmid 1989; T. J. Harries & I. D. Howarth 1996; S.-J. Chang & H.-W. Lee 2020; B.-E. Choi et al. 2020; J.-E. Heo et al. 2021; S.-J. Chang et al. 2023b) can affect the formation of Raman-scattered features.

We will show results of Raman-scattered continuum photons in Section 3.2 and discuss Raman-scattered features of the emission lines of the higher Lyman series ($\text{Ly}\beta$ and $\text{Ly}\gamma$) in Appendix B.

2.3.2 Monte Carlo simulation

The Raman-scattering simulations are performed using the Monte Carlo code STARS (S.-J. Chang & H.-W. Lee 2020), following these steps:

- (i) Define the spherical geometry with uniform HI column density N_{HI} .
- (ii) Generate a flat UV continuum between 940–1060 Å from the central source or Gaussian emission of Lyman series.
- (iii) For each photon, determine the scattering location using the total scattering cross-section (the black solid line in Fig. A1). The details to calculate a free path are described in (iii) of Section 2.2.3.
- (iv) Select the de-excitation channel (i.e. final state of the electron) based on branching ratios computed from the Kramers–Heisenberg formula. If the random number r from 0 to 1 is smaller than the branching ratio to the $n > 1$ state, the photon Raman scatters and converts to an optical or IR photon (e.g. $H\alpha$ from $\text{Ly}\beta$, and $H\beta$ and $\text{Pa}\alpha$ from $\text{Ly}\gamma$). Otherwise, the photon undergoes Rayleigh scattering and remains a UV photon.
- (v) Compute the scattered wavelength λ_{sc} using equation (6), and generate the new photon direction following the Thomson-scattering phase function (equation 9) as Raman scattering is off-resonance scattering (W. J. Karzas & R. Latter 1961). If the photon Raman scatters, it is considered to escape the grid immediately.
- (vi) Repeat steps (iii)–(v) until the photon escapes the grid or undergoes Raman scattering.
- (vii) Back to a step (ii) until the total number of intrinsic photons becomes 10^8 .

The detailed Monte Carlo method of STARS is described in section 2 of S.-J. Chang & H.-W. Lee (2020). We will investigate the formation of the broad wings around hydrogen lines by Raman scattering and estimate their widths in Section 3.2.

2.4 Thomson scattering

Thomson scattering is particularly relevant in the presence of ionized gas in LRDs. The exponential wings observed in many Balmer profiles suggest the influence of electron scattering, as proposed in

previous work (A. Laor 2006; V. Rusakov et al. 2025). Importantly, Thomson scattering produces symmetric broadening across all lines when the emission region is embedded within the scattering medium. Unlike resonance or Raman processes, Thomson scattering can introduce significant line broadening without requiring large gas velocities. This poses a challenge for interpreting Balmer line widths as purely virial in origin, motivating detailed modelling of this mechanism in the LRD context.

2.4.1 Physical mechanism

Thomson scattering is an elastic process in which photons scatter with free electrons, without any change in photon energy in the electron’s rest frame. The scattering cross-section is independent of wavelength (G. B. Rybicki & A. P. Lightman 1986) and given by

$$\sigma_{\text{Th}} \approx 6.65 \times 10^{-25} \text{ cm}^2. \quad (8)$$

The angular distribution of scattered photons is described by the Thomson phase function:

$$P(\theta) = \frac{3}{16\pi} (1 + \cos^2 \theta), \quad (9)$$

where θ is the scattering angle between the incident and scattered directions. With this scattering phase function, the probability of forward and backward scattering is greater than that of scattering in the perpendicular direction.

The broadening of the scattered photon spectrum arises from the thermal motion of electrons. The thermal speed of electrons is

$$v_{\text{th,e}} = \sqrt{\frac{2kT_e}{m_e}} \approx 548 \text{ km s}^{-1} \left(\frac{T_e}{10^4 \text{ K}} \right)^{1/2}, \quad (10)$$

where T_e is the electron temperature. Since the scattering cross-section is wavelength-independent, the resulting broad features from Thomson scattering appear around any emission lines from the same emission region with a similar width in velocity space, if these lines are emitted from the same region.

2.4.2 Monte Carlo simulation

Since Thomson scattering is an elastic process in the electron’s rest frame, like resonance scattering, the Thomson-scattering simulations are performed using the Monte Carlo code RT-SCAT (S.-J. Chang & M. Gronke 2024) with the wavelength-independent cross-section, following these steps:

- (i) Define the spherical geometry with uniform electron density n_e , electron temperature T_e , and compute the Thomson optical depth $\tau_{\text{Th}} = \sigma_{\text{Th}} n_e R$.
- (ii) Generate isotropic photons from the central source, with an intrinsic Gaussian width σ_{src} . Due to the wavelength independence of the Thomson scattering cross-section, the generated photon represents any emission from the central source (e.g. $H\alpha$ and $H\beta$).
- (iii) For each photon, determine the scattering location using the Thomson cross-section σ_{Th} in equation (8) [details are described in (iii) of Section 2.2.3].
- (iv) Compute the new photon direction using the Thomson phase function. The scattered wavelength is identical to the incident one in the rest frame of the electron. However, the scattered wavelengths vary in the observed frame. With a 3D velocity from a Maxwell–Boltzmann distribution $\mathbf{v}_e = v_{\text{th,e}} \hat{\mathbf{v}}$ at an electron temperature T_e , a scattered wavelength in the observed frame λ_s is given by

$$\lambda_s \approx \lambda_i \left[1 + \frac{v_{\text{th,e}}}{c} \hat{\mathbf{v}} \cdot (\hat{\mathbf{k}}_i - \hat{\mathbf{k}}_s) \right], \quad (11)$$

where \hat{k}_i and \hat{k}_s are unit vectors of incident and scattered photons, and λ_i is an incident wavelength in the observed frame; the range of $\hat{v} \cdot (\hat{k}_i - \hat{k}_s)$ term is $[-1, 1]$.

(v) Repeat steps (iii)–(iv) until the photon escapes the grid.

(vi) Back to a step (ii) until the total number of intrinsic photons becomes 10^8 .

The detailed Monte Carlo method of RT-SCAT is described in section 2.5 of S.-J. Chang & M. Gronke (2024; also see A. Laor 2006; H. J. Kim, H.-W. Lee & S. Kang 2007; S.-J. Chang et al. 2018b; H. Y. Choe & H.-W. Lee 2023). We will explore Thomson-scattered features for various electron temperatures T_e and optical depths τ_{Th} , and investigate which environment forms an exponential profile.

3 RESULTS

In this section, we present the simulated spectral features of hydrogen emission lines arising from resonance, Raman, and Thomson scattering processes. Our RT models predict how each mechanism imprints distinct signatures on Balmer lines, such as H α and H β . We systematically explore the effects of key physical parameters for each scattering process. In Section 3.1, we investigate the multitransition effects in the $n = 2$ and 3 states involved in H β scattering and the formation of H α and H β spectra. In Section 3.3, we explore the formation of Thomson-scattered features for various electron temperatures and optical depths. In Section 3.2, we investigate the broad wings around hydrogen emission lines through Raman scattering of UV radiation near the Lyman series.

3.1 Resonant scattering of Balmer lines

3.1.1 Pa α and H α from H β scattering

While H β photons undergo scattering, they can convert to Pa α and H α photons through transitions from the $4p$ state to the $3s$ and $3d$ states, with a branching ratio BR3 of 0.261 (see Fig. 2 and Table 3). To quantify this process, we adopt a monochromatic source emitting H β photons at the line centre and estimate the conversion rate of H β to Pa α via scattering to $n = 3$ states, defined as

$$C_{\text{H}\beta \rightarrow \text{Pa}\alpha} = \frac{\Phi_{\text{H}\beta \rightarrow \text{Pa}\alpha}}{\Phi_{\text{H}\beta, \text{in}}}, \quad (12)$$

where $\Phi_{\text{H}\beta, \text{in}}$ is the number of intrinsic H β photons, and $\Phi_{\text{H}\beta \rightarrow \text{Pa}\alpha}$ is the number of escaping Pa α photons generated by H β scattering. Fig. 4 shows $C_{\text{H}\beta \rightarrow \text{Pa}\alpha}$ as a function of $N_{\text{HI}, 2s}$.

$C_{\text{H}\beta \rightarrow \text{Pa}\alpha}$ increases with increasing $N_{\text{HI}, 2s}$ in Fig. 4 because higher optical depths lead to more scatterings, providing more opportunities for H β photons to convert to Pa α and H α via the $4p \rightarrow 3s/3d$ transition. To understand this trend, we derive an analytic solution for $C_{\text{H}\beta \rightarrow \text{Pa}\alpha}$.

The effective optical depth for the $n = 3$ transition in a spherical geometry is BR3 $\tau_{0, \text{H}\beta}$, where BR3 is the branching ratio to the $n = 3$ state, ~ 0.26 . The number of escaping H β photons consists of those that either escape directly or undergo scattering into the $2s$ state. In other words, they are photons that escape without converting via the $n = 3$ transition. The number of escaping H β photons is thus

$$\Phi_{\text{H}\beta} = \Phi_{\text{H}\beta, \text{in}} e^{-\text{BR3} \tau_{0, \text{H}\beta}}. \quad (13)$$

The number of Pa α photons produced via scattering is $\Phi_{\text{in}} - \Phi_{\text{H}\beta}$, i.e.

$$\Phi_{\text{H}\beta \rightarrow \text{Pa}\alpha} = \Phi_{\text{H}\beta, \text{in}} (1 - e^{-\text{BR3} \tau_{0, \text{H}\beta}}). \quad (14)$$

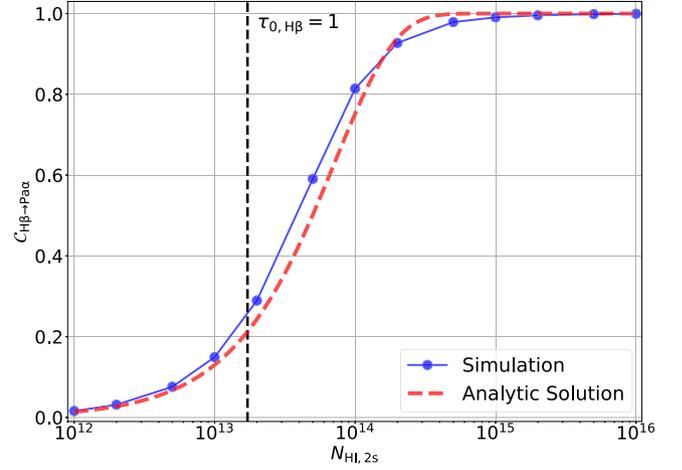


Figure 4. Conversion rate from H β to Pa α through inelastic H β scattering from $n = 2 \rightarrow 4 \rightarrow 3$, $C_{\text{H}\beta \rightarrow \text{Pa}\alpha}$, as a function of $N_{\text{HI}, 2s}$. The blue dots show the simulation results for the monochromatic case, while the red dashed line represents the analytic solution derived in equation (15). The vertical dashed line represents the $N_{\text{HI}, 2s}$ corresponding to an optical depth of unity at the H β line centre ($\tau_{0, \text{H}\beta} = 1$). The conversion rate $C_{\text{H}\beta \rightarrow \text{Pa}\alpha}$ increases with $N_{\text{HI}, 2s}$, as higher optical depths lead to more scatterings and chances of branching via the $4p \rightarrow n = 3$ ($3s/3d$) transition.

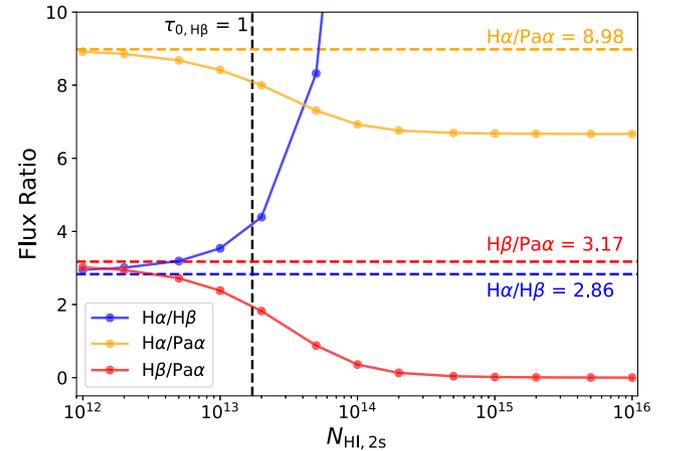


Figure 5. Emergent flux ratios experiencing H β scattering of H α /H β (blue), H β /Pa α (red), and H α /Pa α (orange) as functions of $N_{\text{HI}, 2s}$. The ratio variations originate from H β scattering, as shown in Fig. 4. The vertical dashed line marks the $N_{\text{HI}, 2s}$ corresponding to an optical depth of unity at the H β line centre ($\tau_{0, \text{H}\beta} = 1$). The horizontal dashed lines indicate the line ratios expected under Case B recombination.

Thus, the conversion rate is

$$C_{\text{H}\beta \rightarrow \text{Pa}\alpha} = \frac{\Phi_{\text{H}\beta \rightarrow \text{Pa}\alpha}}{\Phi_{\text{H}\beta, \text{in}}} = 1 - e^{-\text{BR3} \tau_{0, \text{H}\beta}} \quad (15)$$

Fig. 4 shows that the analytic solution and the simulation results match well. For $\tau_{0, \text{H}\beta} > 0.5$, the simulation slightly exceeds the analytic estimate due to multiple scatterings, which increase the effective path length via random walks.

Fig. 5 presents the variation of line flux ratios – H α /H β , H β /Pa α , and H α /Pa α – as functions of $N_{\text{HI}, 2s}$, assuming intrinsic line ratios consistent with Case B recombination. As H β photons begin to convert into Pa α and H α at $N_{\text{HI}, 2s} \gtrsim 10^{13} \text{ cm}^{-2}$, the H α /H β ratio

increases with $N_{\text{HI},2s}$. This trend is similar to the common Balmer decrement due to the dust extinction, but it originates from $\text{H}\beta$ scattering. However, note that the intrinsic spectra used in Fig. 5 are δ -functions at the respective line centres. If, on the other hand, the incident radiation has a broad intrinsic width and photons can escape more easily directly from the central source without significant interaction, the variation in line ratios can be significantly smaller.

3.1.2 Formation of the $\text{H}\alpha$ spectrum

In this section, we examine the emergent $\text{H}\alpha$ spectra for a Gaussian input profile with an intrinsic width of $\sigma_{\text{Src}} = 200 \text{ km s}^{-1}$. Considering that $\text{H}\alpha$ RT involves only the $2s - 3p$ transition – analogous to $\text{Ly}\alpha$ RT with the $1s - 2p$ transition – we anticipate spectral behaviour similar to that of $\text{Ly}\alpha$ – however, typically at much lower optical depths. Fig. 6 illustrates $\text{H}\alpha$ spectra for varying $N_{\text{HI},2s}$, σ_{R} , and v_{exp} using an intrinsic spectral width of $\sigma_{\text{Src}} = 200 \text{ km s}^{-1}$.

In the left panel of Fig. 6, as $N_{\text{HI},2s}$ increases, the double-peaked profiles near the line centre become more pronounced, and the separation between the peaks widens due to an increased number of scatterings. The central panel shows that increasing the random motion of the scattering medium σ_{R} leads to broader double peaks, reflecting enhanced frequency diffusion driven by greater random gas motions.

In the right panel of Fig. 6, an outflow velocity smaller than σ_{Src} (i.e. $v_{\text{exp}} = 50$ and 100 km s^{-1}) shows a P-Cygni profile composed of an absorption-like feature blueward of line centre and a red peak. For larger outflow velocities ($v_{\text{exp}} \geq 500 \text{ km s}^{-1}$), scattering becomes negligible because the medium effectively becomes optically thin for the intrinsic photons.

In summary, the spectral behaviour of $\text{H}\alpha$ is similar to that of $\text{Ly}\alpha$, albeit $\text{Ly}\alpha$ is typically studied at much larger column densities $N_{\text{HI}} \gtrsim 10^{17} \text{ cm}^{-2}$. Leaving aside the doublet structure, the emergent $\text{H}\alpha$ spectrum resembles the one typically found for metal resonance lines observed in outflowing environments (e.g. Mg II , J. X. Prochaska, D. Kasen & K. Rubin 2011; S.-J. Chang & M. Gronke 2024).

3.1.3 $\text{H}\beta$ spectrum

In this section, we consider the intrinsic $\text{H}\beta$ emission from a central source, assuming a width of $\sigma_{\text{Src}} = 200 \text{ km s}^{-1}$. The RT of $\text{H}\beta$ involves transitions from the $4p$ state to both the $2s$ state and the $3s/3d$ states. The scattering characteristics of $\text{H}\beta$ differ from those of $\text{H}\alpha$ and $\text{Ly}\alpha$ because of the transition to the $3s/3d$ states, as discussed in Section 3.1.1.

Fig. 7 shows $\text{H}\beta$ spectra for various $N_{\text{HI},2s}$, σ_{R} , and v_{exp} . In the left and central panels for a static medium ($v_{\text{exp}} = 0 \text{ km s}^{-1}$), the spectra do not exhibit clear double-peaked features like those observed in $\text{H}\alpha$ spectra in Fig. 6. Instead, the profiles look like simple Voigt absorption profiles centred at the intrinsic wavelength. This occurs because most $\text{H}\beta$ photons rarely experience sufficient scattering events to form a double-peaked profile before converting into $\text{Pa}\alpha$. As discussed in Section 3.1.1, when the optical depth at $\text{H}\beta$ exceeds ~ 10 , most $\text{H}\beta$ photons undergo scattering through the $4p \rightarrow 3s/3d$ transition, converting into $\text{Pa}\alpha$ and $\text{H}\alpha$ photons (cf. Fig. 2). The branching ratio to $n = 3$ states ~ 0.26 is non-negligible, implying that it is statistically unlikely for $\text{H}\beta$ photons to remain in the $\text{H}\beta$ state after more than about four scatterings ($\sim 1/\text{BR}_3$).

In the right panel of Fig. 7, for the slow outflow case smaller than $\sigma_{\text{Src}} = 200 \text{ km s}^{-1}$ ($v_{\text{exp}} = 50$ and 100 km s^{-1}), a red peak emerges due to the scattering in the outflowing medium. This behaviour contrasts with the static case because the frequency shifts

in an outflow allow some scattered $\text{H}\beta$ photons to escape before conversion to $\text{Pa}\alpha$. Nevertheless, since conversion to $\text{Pa}\alpha$ is still efficient, the red peak in $\text{H}\beta$ remains weaker compared to that seen in the corresponding $\text{H}\alpha$ spectra in the right panel of Fig. 6. We will discuss the differences between $\text{H}\alpha$ and $\text{H}\beta$ spectra in the following section.

3.1.4 Comparing $\text{H}\alpha$, $\text{H}\beta$, and $\text{Pa}\alpha$ spectra

Fig. 8 illustrates the impact of multiple-transition scattering of $\text{H}\beta$ photons on the emergent line profiles of $\text{H}\beta$, $\text{H}\alpha$, and $\text{Pa}\alpha$, assuming only $\text{H}\beta$ emission as input radiation. In the top panel for the static medium ($v_{\text{exp}} = 0 \text{ km s}^{-1}$), the $\text{H}\beta$ spectrum exhibits a central absorption-like feature without clear double peaks, indicating significant photon conversion into $\text{Pa}\alpha$ via the $4p \rightarrow n = 3$ transition. The $\text{Pa}\alpha$ spectrum emerges as a symmetric Gaussian profile, shaped by the random motion σ_{R} of the medium. Simultaneously, this conversion generates additional $\text{H}\alpha$ photons, which are emitted at the line-centre wavelength and subsequently scatter, forming a double-peaked profile.

In the bottom panel of Fig. 8, the presence of the outflow ($v_{\text{exp}} = 50 \text{ km s}^{-1}$) significantly modifies the spectral shape. The $\text{H}\beta$ line exhibits a prominent P-Cygni profile. The partial photons on the blue side of the line centre are scattered into the $n = 3$ states due to the outflowing medium, converting into $\text{Pa}\alpha$ photons that escape as a blueshifted emission. The simultaneously generated $\text{H}\alpha$ photons continue to scatter, forming a redshifted peak. The effect of $\text{H}\beta$ scattering in the outflowing medium creates distinct observational signatures, blueshifted $\text{Pa}\alpha$ and redshifted $\text{H}\alpha$ emission features.

To investigate the observational impact of $\text{H}\beta$ scattering, we assume the central source emits intrinsic Gaussian profiles for $\text{H}\alpha$, $\text{H}\beta$, and $\text{Pa}\alpha$ with intrinsic width $\sigma_{\text{Src}} = 200 \text{ km s}^{-1}$ and relative flux ratios following Case B recombination. Since the $n = 3$ population of neutral hydrogen is negligible due to its short lifetime, intrinsic $\text{Pa}\alpha$ photons escape directly without scattering, resulting in a symmetric Gaussian profile.

Fig. 9 presents the observable emergent spectra (i.e. considering both the intrinsic as well as converted radiation) of $\text{H}\alpha$, $\text{H}\beta$, and $\text{Pa}\alpha$ in the outflowing case. Both $\text{H}\alpha$ and $\text{H}\beta$ exhibit P-Cygni profiles, with the red peak of $\text{H}\alpha$ stronger than that of $\text{H}\beta$, due to additional photons produced via the $n = 3$ transition. The $\text{Pa}\alpha$ profile has an enhanced blueward peak due to the direct conversion from $\text{H}\beta$ scattering, as shown in Fig. 8.

In summary, even though the optical depths of $\text{H}\alpha$ and $\text{H}\beta$ are of similar magnitudes due to the same $N_{\text{HI},2s}$ column density and comparable cross-sections (equation 3), their RT processes differ substantially. While $\text{H}\alpha$ RT closely mirrors that of $\text{Ly}\alpha$ in a low optical depth regime (i.e. dominated by resonance scattering without line conversion), $\text{H}\beta$ photons cannot undergo a large number of scatterings due to the large branching ratio into the $n = 3$ state. As a result, the $\text{H}\alpha$ line can be significantly reshaped due to resonant RT, and $\text{H}\beta$ only shows absorption-like features (with additional $\text{Pa}\alpha$ emergent at the corresponding velocity offset). We will discuss the potential impact of resonant scattering on observed Balmer line profiles in Sections 4.2 and 4.4.

3.2 Raman scattering

Fig. 10 shows the Raman-scattered features around various hydrogen lines, $\text{H}\alpha$, $\text{H}\beta$, $\text{H}\gamma$, $\text{Pa}\alpha$, $\text{Pa}\beta$, and $\text{Br}\alpha$ from the UV continuum near $\text{Ly}\beta$, $\text{Ly}\gamma$, and $\text{Ly}\delta$. As N_{HI} increases, Raman features get broader, reflecting the fact that a wider range of UV continuum photons near

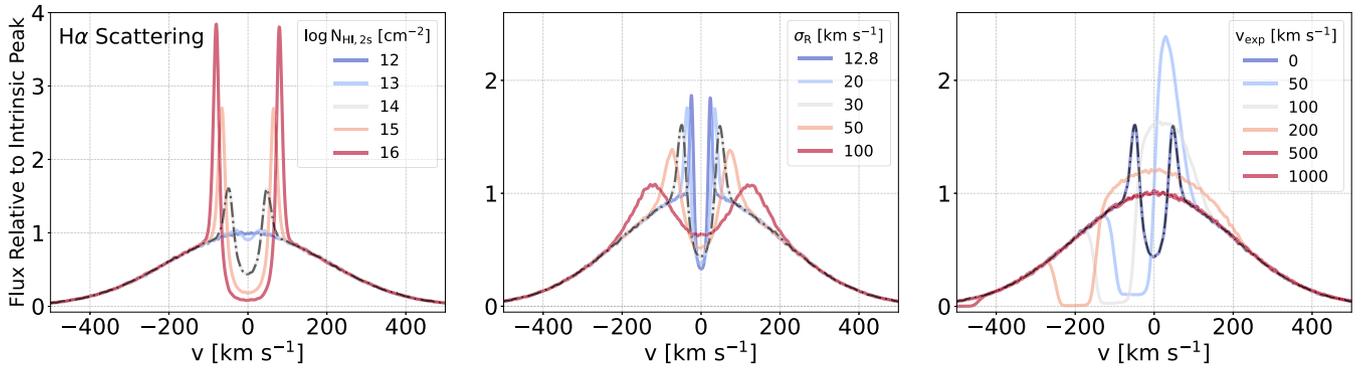


Figure 6. Spectra of resonance-scattered H α for different values of $N_{\text{HI},2s}$ (left panel), random velocities σ_R (centre panel), and bulk outflow speeds v_{exp} (right panel). The spectra are normalized by dividing the peak flux of intrinsic radiation; thus, the y-axis is the relative flux. Fiducial parameters are set to $N_{\text{HI},2s} = 10^{14} \text{ cm}^{-2}$, $v_{\text{exp}} = 0 \text{ km s}^{-1}$, and $\sigma_R = 30 \text{ km s}^{-1}$. The intrinsic Gaussian emission profile width σ_{src} is fixed at 200 km s^{-1} . Spectral variations closely resemble those of Ly α under similar optical depth conditions, although typical Ly α optical depths are orders of magnitude larger.

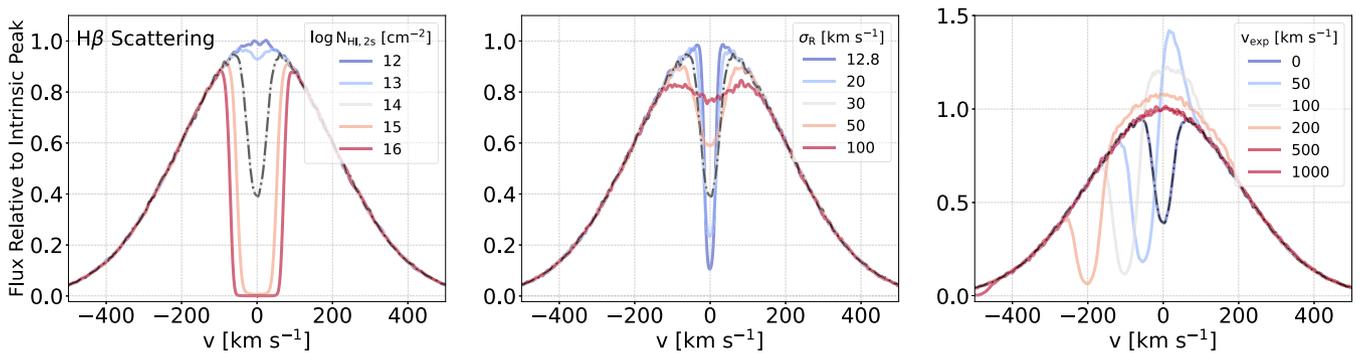


Figure 7. Spectra of resonance-scattered H β for various $N_{\text{HI},2s}$ (left), σ_R (centre), and v_{exp} (right). The parameters of the fiducial case (black dot–dashed line) and the range of parameters for each panel are the same as those in Fig. 6. The spectra are normalized by dividing the peak of intrinsic radiation at $v = 0 \text{ km s}^{-1}$; y-axis represents the relative flux. Unlike the double-peaked structures prominently visible in the H α spectra in the left and centre panels of Fig. 6, the H β line generally shows a central absorption-like feature due to significant photon conversion to Pa α via scattering into $n = 3$ states.

the Lyman series undergo Raman scattering to form broad features around hydrogen lines (S.-J. Chang et al. 2015, 2018b).

For example, UV continuum near Ly β produces broad wings around H α via Raman scattering into the $n = 2$ state. Similarly, Ly γ photons scatter into both $n = 2$ and 3 states, generating Raman features near H α and Pa α , respectively. Ly δ photons also contribute via Raman scattering into the $n = 4$ state, producing Br α features. At high $N_{\text{HI}} (> 10^{21} \text{ cm}^{-2})$, the incident flat continuum is fully imprinted onto the Raman-scattered features, resulting in a flattened spectral shape near each line centre. The slight slopes in these flat regions originate from the energy conservation inherent in the Raman-scattering process (equation 7).

The left panel of Fig. 10 shows that the Raman H α feature is broader than those of H β and H γ at the same N_{HI} . This trend is driven by two factors: (1) the larger Raman-scattering cross-section near Ly β compared to Ly γ and Ly δ (see Fig. A1), and (2) the different wavelength-broadening factors in equation (7). As a result, the α line of each hydrogen series (Balmer, Paschen, and Brackett) exhibits the broadest Raman wings relative to the higher order transitions, assuming a flat input continuum. These spectral trends offer a potential observational diagnostic: broader wings of Balmer lines (e.g. H α compared to H β) may indicate a Raman-scattered wing. We discuss these observational signatures of Raman scattering further in Section 4.3.

We further compare the widths and profile shapes of Raman-scattered features around hydrogen lines. In Fig. 11, when $N_{\text{HI}} = 10^{20-23} \text{ cm}^{-2}$, the width of Raman H α features spans ~ 1000 to $30\,000 \text{ km s}^{-1}$, scaling approximately as $N_{\text{HI}}^{1/2}$. Moreover, the width varies across hydrogen lines depending on their origin of Lyman series; H α , H β , and H γ arise from Raman scattering of Ly β , Ly γ , and Ly δ , respectively. Because Ly β has both a higher scattering cross-section and a larger Raman wavelength shift (see Fig. A1 and equation 7), H α wings are significantly broader – typically by a factor of ~ 3 – compared to H β .

In summary, Raman scattering produces broader features at higher N_{HI} due to increased optical depth and interaction range with the UV continuum. Continuum Raman-scattered features show the different widths of each hydrogen line. If H α wings are significantly broader than H β , it suggests a contribution from Raman scattering of the UV continuum. However, this width contrast is not significant when Ly β and Ly γ appear as intrinsic emission features in the input radiation, rather than a flat continuum. We discuss this case in more detail in Section 4.3 and Appendix B.

3.3 Thomson scattering

Fig. 12 shows simulated results of Thomson-scattered line profiles for various electron temperatures $T_e = 10^{4-6} \text{ K}$ and Thomson optical

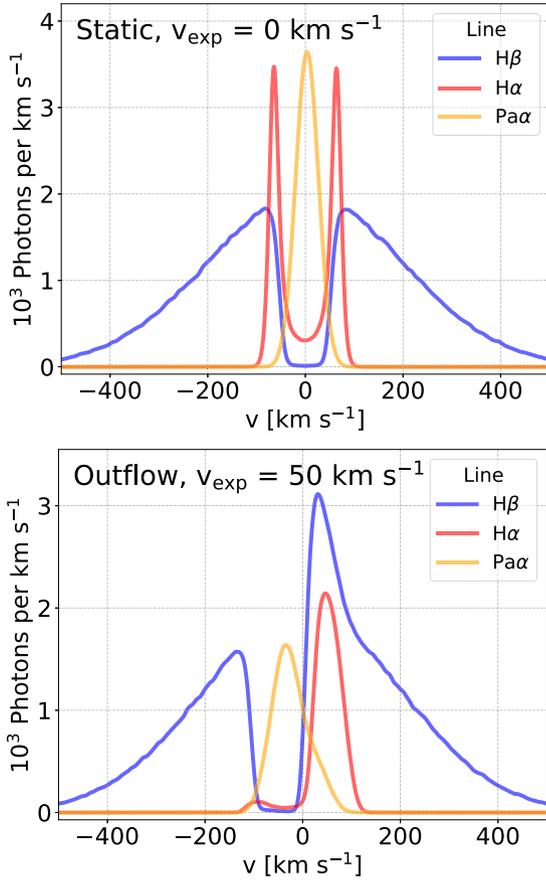


Figure 8. Simulated spectra as a consequence of resonantly scattered H β RT, including $4p \rightarrow n=2$ and $4p \rightarrow n=3$ transitions and assuming only H β Gaussian emission with $\sigma_{\text{src}} = 200 \text{ km s}^{-1}$. The colours of solid lines represent simulated spectra of H β (blue), H α (red), and Pa α . H α and Pa α spectra are composed of photons produced by H β scattering to the $n=3$ state, while H β spectra are composed of scattered and directly escaping H β . y-axis is the number of photons per km s^{-1} to show how many photons convert from H β to Pa α and H α . $N_{\text{HI},2s}$ and σ_{R} are fixed at 10^{15} cm^{-2} and 30 km s^{-1} , respectively. The top panel shows spectra for a static medium ($v_{\text{exp}} = 0 \text{ km s}^{-1}$), while the bottom panel presents results for an outflowing medium ($v_{\text{exp}} = 50 \text{ km s}^{-1}$).

depth $\tau_{\text{Th}} = 0.1 - 10$. Since the dependence on the input radiation is negligible if $\sigma_{\text{src}} < v_{e,\text{th}}$ (i.e. in most cases), we assume a width of the intrinsic emission $\sigma_{\text{src}} = 50 \text{ km s}^{-1}$ to focus on the broad wing features due to Thomson scattering. In the left panel of Fig. 12, the width of the scattered wings increases with increasing T_e , following the trend in electron thermal velocity $v_{\text{th},e}$ in equation (10). Since Thomson scattering is independent of photon wavelength, the resulting broad wings appear around all emission lines with similar shapes, as discussed in Section 2.4.

The bottom panel of Fig. 12 shows the dependence on the optical depth τ_{Th} for a fixed electron temperature $T_e = 10^4 \text{ K}$. In the optically thin case ($\tau_{\text{Th}} \leq 1$), where single scattering dominates, the profiles exhibit exponential wings, consistent with the analytical form $\propto \exp(-v/v_0)$ (R. A. Sunyaev & L. G. Titarchuk 1980; H.-W. Lee 1999; A. Laor 2006). With the single scattering approximation, the analytic solution of the Thomson-scattered profile is given by

$$f(v) = f_0 e^{-v/v_0}, \quad (16)$$

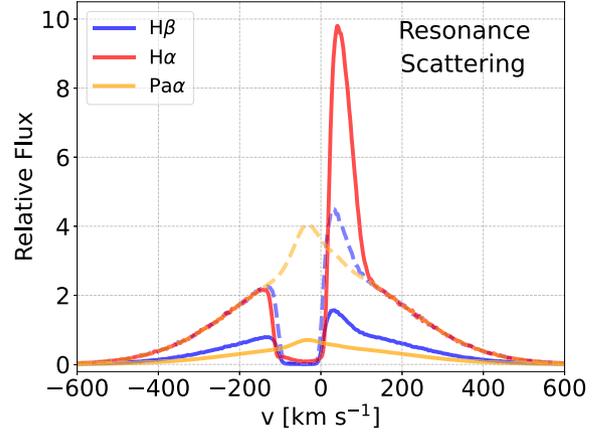


Figure 9. Emergent spectra of H α , H β , and Pa α after RT of H α and H β resonance scattering through an outflow medium with $v_{\text{exp}} = 50 \text{ km s}^{-1}$ and $N_{\text{HI},2s} = 10^{15} \text{ cm}^{-2}$. The widths of intrinsic profiles of H α , H β , and Pa α , σ_{src} , are fixed at 200 km s^{-1} , assuming the flux ratios between lines under Case B recombination. The blue solid line is identical to the H β spectrum in the bottom panel of Fig. 8. The red (orange) solid line shows the integrated H α (Pa α) spectrum combining the emergent spectrum of intrinsic emission and those produced via H β scattering to the $n=3$ state in bottom panel of Fig. 8. While the *solid lines* show the correct (observable) flux ratios, the *dashed curves* are normalized to the wings, illustrating their relative spectral shapes.

where $v_0 \approx 1.1 v_{e,\text{Th}}$ (A. Laor 2006). This analytic solution (black dashed line) is akin to the Thomson wing at $\tau_{\text{Th}} = 0.1$, in the bottom panel of Fig. 12. When τ_{Th} increases, the profile gets broader as a higher optical depth provides the opportunity for additional scattering. However, as still single scattering mainly determines the formation of Thomson-scattered features, the profile follows the exponential profile well at $\tau_{\text{Th}} \leq 1$.

In the optically thick case ($\tau_{\text{Th}} \gtrsim 2$), multiple scatterings modify the profile shape, deviating from the exponential form as can be seen in the bottom panel of Fig. 12. This results in a mismatch between the simulated profiles (solid lines) and the exponential fits (dashed lines). The cumulative effects of repeated velocity shifts and directional variations through multiple scatterings result in a Gaussian-like profile near the line centre with a power-law wing (R. A. Sunyaev & L. G. Titarchuk 1980; L. Dessart et al. 2009).

To quantify the broadening, we estimate the full width at half-maximum (FWHM) of the Thomson-scattered profiles. In the top panel Fig. 13, we show the FWHM as a function of electron temperature. As expected from equations (16) and (10), the FWHM increases as $T_e^{1/2}$, independent of τ_{Th} . In the bottom panel, we show how the spectral broadening depends on the Thomson optical depth. FWHM remains constant for $\tau_{\text{Th}} < 1$, where single scattering dominates, but increases approximately linearly with τ_{Th} in the optically thick regime due to the cumulative effect of multiple scatterings.

In summary, Thomson scattering produces symmetric, broad line wings whose width scales with $T_e^{1/2}$ and τ_{Th} . In the optically thin regime, the line profile exhibits exponential profiles – a key diagnostic feature that has been identified in broad Balmer emission of some AGNs and LRDs (e.g. A. Laor 2006; V. Rusakov et al. 2025). However, in the optically thick case ($\tau_{\text{Th}} \geq 5$), the Thomson-scattered feature becomes non-exponential (as shown in the bottom panel of Fig. 12). Moreover, if the Thomson-scattering medium is outflowing, bulk motions can cause asymmetry in the broad scattered wings, as demonstrated in Appendix C (see also L. H. Auer & D. van

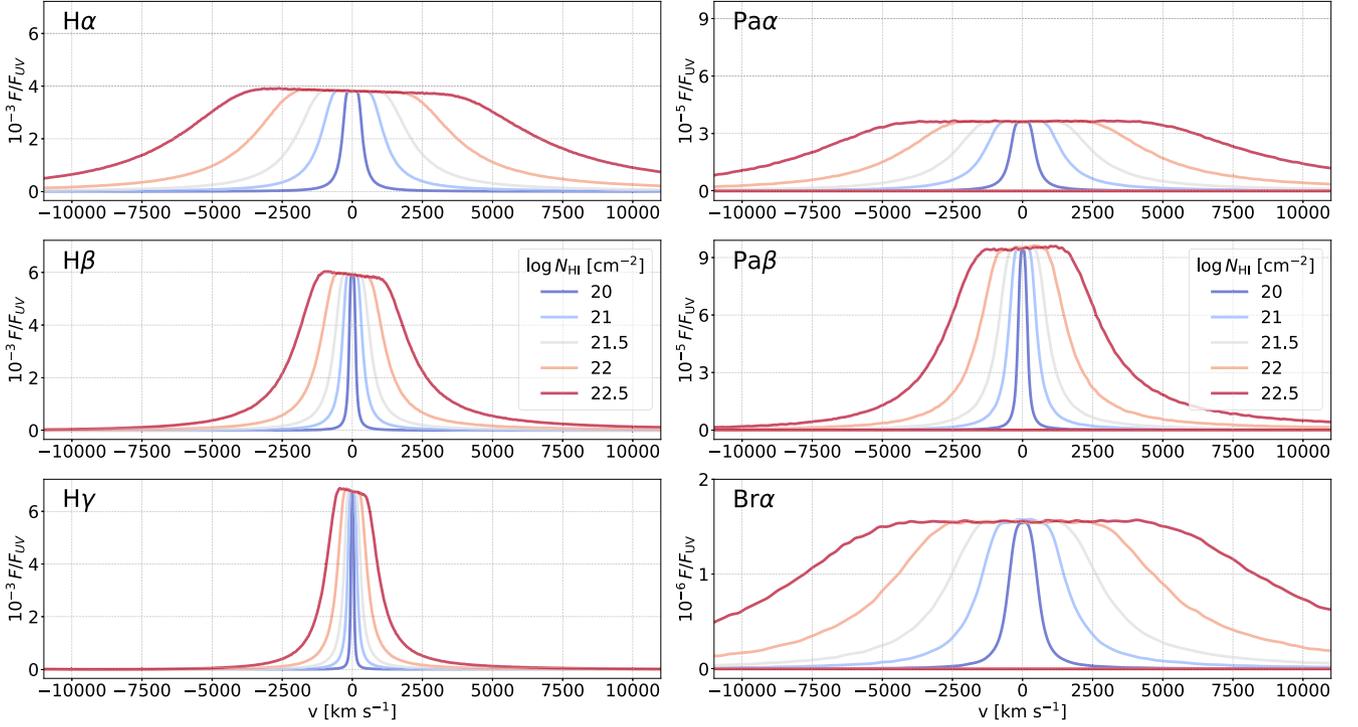


Figure 10. Raman-scattered features of hydrogen lines for various H I column densities, $N_{\text{HI}} = 10^{20-22.5} \text{ cm}^{-2}$. The vertical axis F/F_{UV} denotes the flux of the Raman feature normalized by the incident UV continuum near the corresponding Lyman-series line. Specifically, F_{UV} refers to the continuum around Ly β , Ly γ , and Ly δ for H α , H β , and H γ , respectively. Left: Raman features in Balmer lines: H α (top), H β (middle), and H γ (bottom). Right: Raman features in IR transitions: Pa α (top), Pa β (middle), and Br α (bottom). As N_{HI} increases, the wings broaden due to Raman scattering of UV continuum photons over a wider frequency range.

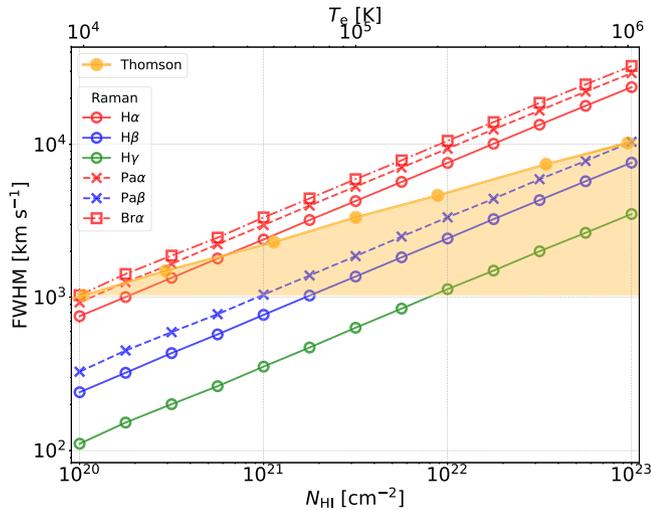


Figure 11. FWHM of Raman-scattered features. FWHM is plotted as a function of $N_{\text{HI}} = 10^{20-23} \text{ cm}^{-2}$. The red, blue, and green solid lines represent Raman features of Balmer lines H α , H β , and H γ , respectively, while dashed and dot-dashed lines correspond to Paschen and Brackett lines. The orange solid line represents FWHM of Thomson-scattered features as a function of T_e at fixed optical depth $\tau_{\text{Th}} = 1$. Unlike Thomson scattering – which yields the same width for all hydrogen lines due to its wavelength-independent cross-section – Raman scattering produces broader H α wings than H β or H γ , driven by differences in scattering cross-section and Raman wavelength shift.

Blerkom 1972). Observationally, the absence of a strong red–blue asymmetry in LRD Balmer wings would constrain $v_{\text{exp}}/v_{\text{th,e}} \ll 1$ in the H II scattering layer. We will further discuss the implications of a Thomson origin for Balmer line broadening in Section 4.

4 DISCUSSION

4.1 Observational signatures of scattering mechanisms

We have investigated three distinct RT processes that can shape the observed profiles of hydrogen emission lines in LRDs:

(i) Resonance scattering of Balmer lines with hydrogen atoms in the $2s$ state can produce double-peaked or P-Cygni-like features, influenced by the kinematics of the medium (i.e. v_{exp} and σ_{R} in the geometry used here). Furthermore, resonance scattering of H β affects the flux ratio of hydrogen lines, as multiple H β scatterings convert H β photons to Pa α and H α via de-excitation to the $n = 3$ state (see Fig. 4 and Section 3.1.1).

(ii) Raman scattering of UV continuum photons near the Lyman series by ground-state hydrogen atoms can generate broad wings around Balmer lines, with widths up to several tens of thousands of km s^{-1} , depending primarily on the H I column density N_{HI} (with the width scaling as $N_{\text{HI}}^{1/2}$; see Section 3.2 and specifically Fig. 11).

(iii) Thomson scattering of Balmer line photons by free electrons leads to symmetric broadening, with widths proportional to the electron’s thermal speed $v_{\text{th,e}}$. While mostly this leads to exponential wings ($\propto \exp(-v/v_{\text{th,e}})$), in the optically thin case (Thomson optical depth $\tau_{\text{Th}} < 1$), Thomson-scattered features exhibit potentially observable deviations (cf. Section 3.3 and Fig. 12).

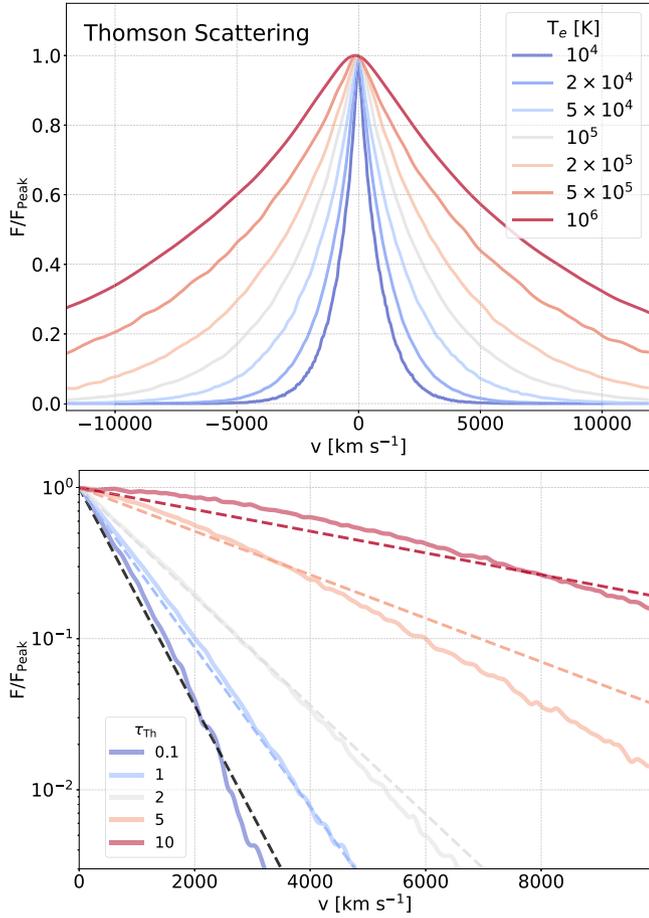


Figure 12. Spectra composed of Thomson-scattered photons, normalized by the peak flux at the line centre. Top: Thomson-scattered profiles for various electron temperatures $T_e = 10^4\text{--}6$ K at fixed optical depth $\tau_{\text{Th}} = 1$, plotted in linear scale. The width of the line wings increases with T_e , corresponding to higher electron thermal speeds. Bottom: Thomson-scattered profiles in the red wing of the line centre at fixed $T_e = 10^4$ K for various optical depths $\tau_{\text{Th}} = 0.1\text{--}10$, plotted in logarithmic scale. Solid lines show Monte Carlo simulation results, and dashed lines represent exponential fits. At low optical depths ($\tau_{\text{Th}} < 1$), profiles follow an exponential shape, while at higher depths ($\tau_{\text{Th}} \gtrsim 2$), multiple scatterings distort the profile from the exponential form.

Fig. 11 shows a summary of the emergent line width due to Raman and Thomson scattering. Importantly, the broad components of hydrogen emission lines arising from Raman scattering exhibit different widths for each line, while Thomson scattering produces similar broadening across all lines, emitted from the same emission region. One should note that if indeed the broad component of LRDs originates from Raman or Thomson scattering, it should be polarized (H.-W. Lee & J.-H. Yun 1998; H.-W. Lee 1999; H. J. Kim et al. 2007) and will be observable through future instruments and telescopes such as the Habitable Worlds Observatory (C. Neiner, A. Girardot & J.-M. Reess 2025). These distinctions provide a diagnostic for identifying the dominant scattering mechanism in LRDs. Tables 4 summarize the key properties of each process.

4.2 Can resonance scattering broaden the observed $H\beta$ profile?

Resonance scattering leads to both frequency and spatial diffusion of photons. For example, resonantly scattered $\text{Ly}\alpha$ emission typically exhibits a double-peaked profile in static media (D. A. Neufeld 1990).

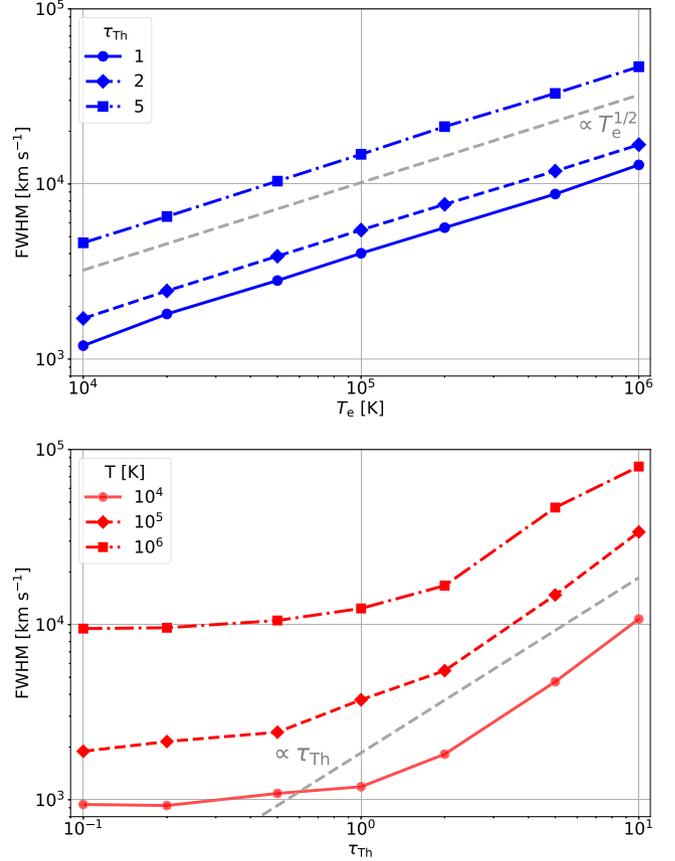


Figure 13. FWHM of Thomson-scattered features. Top: FWHM as a function of electron temperature T_e for different optical depths: $\tau_{\text{Th}} = 1$ (solid), 2 (dashed), and 5 (dot-dashed). The grey dashed line indicates the scaling $\text{FWHM} \propto T_e^{1/2}$, consistent with thermal Doppler broadening. Bottom: FWHM versus τ_{Th} for fixed $T_e = 10^4$ K (solid), 10^5 K (dashed), and 10^6 K (dot-dashed). In the optically thick regime ($\tau_{\text{Th}} > 1$), FWHM increases with τ_{Th} due to multiple scatterings. In the optically thin case ($\tau_{\text{Th}} < 1$), FWHM remains nearly constant and is dominated by single scattering.

The absorption features seen in the Balmer lines of LRDs indicate an enhanced $n = 2$ population (I. Juodžbalis et al. 2024, 2025; J. Matthee et al. 2024; K. Inayoshi & R. Maiolino 2025; R. P. Naidu et al. 2025), suggesting that the gas in these systems is optically thick to Balmer transitions, thus enabling resonance scattering. R. P. Naidu et al. (2025) proposed that the double-peaked $H\beta$ profile in LRDs originates from resonance scattering, analogous to $\text{Ly}\alpha$. However, the presence of multiple branching channels at the $n = 4$ level makes $H\beta$ resonance scattering fundamentally different from that of $\text{Ly}\alpha$. This section explores whether resonance scattering of $H\beta$ in the static medium can realistically account for the observed $H\beta$ features.

For $H\alpha$, the emergent profile in a static medium shows a characteristic double peak due to resonance scattering (see Fig. 6), like $\text{Ly}\alpha$. However, as discussed in Section 3.1.1, in the optically thick regime ($\tau_{H\beta} \gtrsim 10$), most $H\beta$ photons undergo conversion to $\text{P}\alpha$ and $\text{H}\alpha$ via transitions to the $n = 3$ state. As a result, the emergent $H\beta$ spectrum in these conditions primarily reflects photons that escape directly, with an absorption-like dip at the line centre (see the top panel of Fig. 8). Most scattered $H\beta$ photons near the line centre thus escape the medium as $\text{P}\alpha$ or $\text{H}\alpha$ photons.

To test whether $H\beta$ resonance scattering can be a dominant broadening mechanism, we perform RT simulations in a static medium

Table 4. Summary of scattering mechanisms in LRDs.

Process	Medium	Spectral features	Broadening
Resonance ($H\alpha$, $H\beta$)	H I ($n = 2, 2s$)	Double peaks, P-Cygni, variation in the $H\alpha/H\beta$ flux ratio	–
Raman (UV \rightarrow optical/IR)	H I ($n = 1, 1s$)	Broad wings with different widths for each line, up to $\sim 10^4$ km s $^{-1}$	$\propto \sqrt{N_{\text{HI}}}$
Thomson	Free electrons	Symmetric broadening with same width for all lines, up to $\sim 10^3$ km s $^{-1}$	$\propto \sqrt{T_e}, \tau_{\text{Th}}$

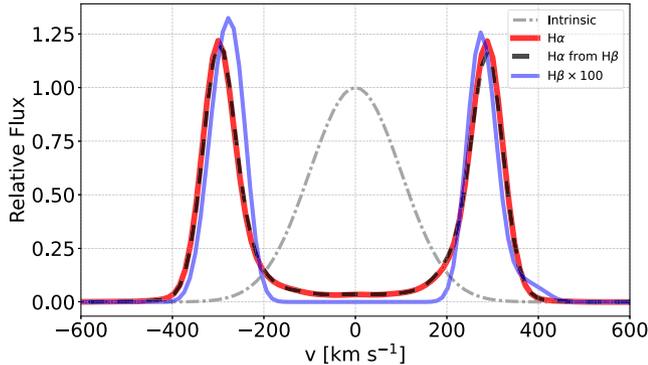


Figure 14. $H\alpha$ and $H\beta$ spectra at $N_{\text{HI}} = 10^{17}$ cm $^{-2}$, $\sigma_{\text{Ran}} = 100$ km s $^{-1}$, and $\sigma_{\text{src}} = 100$ km s $^{-1}$. The intrinsic flux ratio of $H\alpha$ and $H\beta$ is fixed at 1 to compare profiles of emergent spectra. The solid and dashed red lines represent $H\alpha$ spectra from intrinsic $H\alpha$ and $H\beta$, respectively. The blue solid line is $H\beta$ spectra multiplied by a factor of 100. At high N_{HI} and σ_{Ran} , $H\beta$ can be broadened by scattering, but its flux is extremely low as most photons convert to $H\alpha$ through numerous scatterings.

($v_{\text{exp}} = 0$ km s $^{-1}$) with a high column density of $N_{\text{HI},2s} = 10^{17}$ cm $^{-2}$ (i.e. $\tau_{0,H\beta} > 10^3$) and random velocity dispersion $\sigma_{\text{R}} = 100$ km s $^{-1}$. We inject $H\alpha$ and $H\beta$ photons from a central source, each with intrinsic width $\sigma_{\text{src}} = 100$ km s $^{-1}$ and equal flux, ensuring all photons undergo multiple scatterings. Fig. 14 shows the emergent $H\alpha$ spectrum from intrinsic $H\alpha$ (solid red line) and from intrinsic $H\beta$ emission (dashed black line). The two spectra are nearly identical, as most $H\beta$ photons are rapidly converted to $H\alpha$ in the inner scattering region and then escape via subsequent scatterings.

Although the vast majority of $H\beta$ photons are converted to $P\alpha$ and $H\alpha$ (≈ 99 per cent in this case), a small fraction emerges as $H\beta$. The emergent $H\beta$ spectrum (blue line, scaled by a factor of 100 for visibility) in Fig. 14 exhibits a clear double-peaked profile. Notably, the peak separation in $H\beta$ is similar to that in $H\alpha$, since $\tau_{0,H\alpha} \approx \tau_{0,H\beta}$. In summary, $H\beta$ spectra can indeed develop a double-peaked shape via resonance scattering. However, due to the efficient conversion to $P\alpha$, the emergent $H\beta$ flux is several orders of magnitude weaker than $H\alpha$. Since the flux ratio of $H\alpha/H\beta$ in LRDs is not exceeding a hundred, resonance scattering of $H\beta$ is unlikely to be the dominant broadening mechanism.¹

¹Note that if the H I gas is optically thick not only to Balmer lines but also to the Paschen series – due to a large $n = 3$ population – then $P\alpha$ photons may also undergo scattering. In this case, the emergent $H\beta$ flux can increase, and $P\alpha$ shows a double-peaked profile. In gas that is optically thick to both Lyman and Balmer series, the scattering behaviours of $\text{Ly}\beta$ and $H\alpha$ resemble those of $H\beta$ and $P\alpha$ in gas that is optically thick to Balmer and Paschen series, respectively. S.-J. Chang et al. (2018a) explored $H\alpha$ and $\text{Ly}\beta$ emission via resonance scattering and found both can exhibit double-peaked profiles.

4.3 Detection of Raman broad wing

Raman-scattered wings around $H\alpha$ and $H\beta$ have been observed in a variety of astrophysical environments, such as symbiotic stars (H. Nussbaumer et al. 1989; S.-J. Lee & S. Hyung 2018; S.-J. Chang et al. 2018b), planetary nebulae (H.-W. Lee & S. Hyung 2000; H.-W. Lee et al. 2006; L. F. Miranda, J. M. Torrelles & J. Lillo-Box 2022), and H II regions (M. A. Dopita et al. 2016; W. J. Henney 2021). Theoretical studies have also predicted Raman broad wings in AGNs (S.-J. Chang et al. 2015; M. Kokubo & Y. Harikane 2025). The strength of the Raman-scattering wings depends on the input UV radiation. If all incident UV photons are fully Raman scattered, the resulting scattered features have strengths of approximately $\sim 10^{-3}$ relative to the input UV flux (see Fig. 10) due to less energetic scattered photons (equation 5) and Raman broadening (equation 7). Therefore, environments where Raman scattering is observed typically require both strong UV radiation and an optically thick H I medium.

LRDs can provide such conditions, given their strong Balmer breaks and V-shaped SEDs (D. D. Kocevski et al. 2023; J. E. Greene et al. 2024; J. Matthee et al. 2024; H. B. Akins et al. 2025b; A. J. Taylor et al. 2025b). However, the broad wings around $H\alpha$ and $H\beta$ observed in LRDs exhibit similar widths over 1000 km s $^{-1}$ (M. Brazzini et al. 2025; I. Juodžbalis et al. 2025; F. D’Eugenio et al. 2025a, b), which contrasts with the basic Raman-scattering expectation (see Fig. 11) which predicts, e.g. an FWHM ratio of ~ 3 between $H\alpha$ and $H\beta$.

Thus, it seems that Raman scattering is unlikely to be important for LRDs. However, it is important to note that Raman scattering to the $n = 2$ state, specifically the transition $1s \rightarrow np \rightarrow 2s$, leaves an electron in the metastable $2s$ state and can enhance the $n = 2$ population, increasing the optical depth for resonance scattering of Balmer lines. Therefore, although it likely does not shape the emergent spectrum, Raman scattering of continuum photons can still affect the RT dynamics in this manner.

In addition, as shown in Appendix B, Raman scattering of emission lines from the Lyman series can, in fact, produce broad wings of comparable width for both $H\alpha$ and $H\beta$. Assuming the same width of $\text{Ly}\beta$ and $\text{Ly}\gamma$ emissions, the widths of Raman-scattered $H\alpha$ and $H\beta$ are ≈ 6.4 and 5.0 [cf. equation (7) and Section 2.3] times broader than those of the Lyman series, resulting in the width ratio of $H\alpha$ and $H\beta \approx 1.28$. This is because the Raman-scattered profiles reflect the velocity distribution of the source emission rather than that of the Raman cross-section. This implies that for observed widths of ~ 1000 km s $^{-1}$ for LRDs in $H\alpha$, the width of $\text{Ly}\beta$ needs to be ~ 150 km s $^{-1}$, which is well within the observed narrow component of LRDs.

In this case, to form the broad feature with 10 per cent of $H\alpha$ ($H\beta$) flux by Raman scattering, the flux ratio of $\text{Ly}\beta/H\alpha \approx 0.64$ ($\text{Ly}\gamma/H\beta \approx 0.5$) is required, assuming $\text{Ly}\beta$ ($\text{Ly}\gamma$) photons are fully Raman scattered due to high N_{HI} . Although a higher Lyman series emission, such as $\text{Ly}\beta$ and $\text{Ly}\gamma$, is not allowed under Case B recombination, the flux ratios of $\text{Ly}\beta/H\alpha$ and $\text{Ly}\gamma/H\beta$ are ≈ 5 and 2 under Case A recombination (P. J. Storey & D. G. Hummer 1995), respectively,

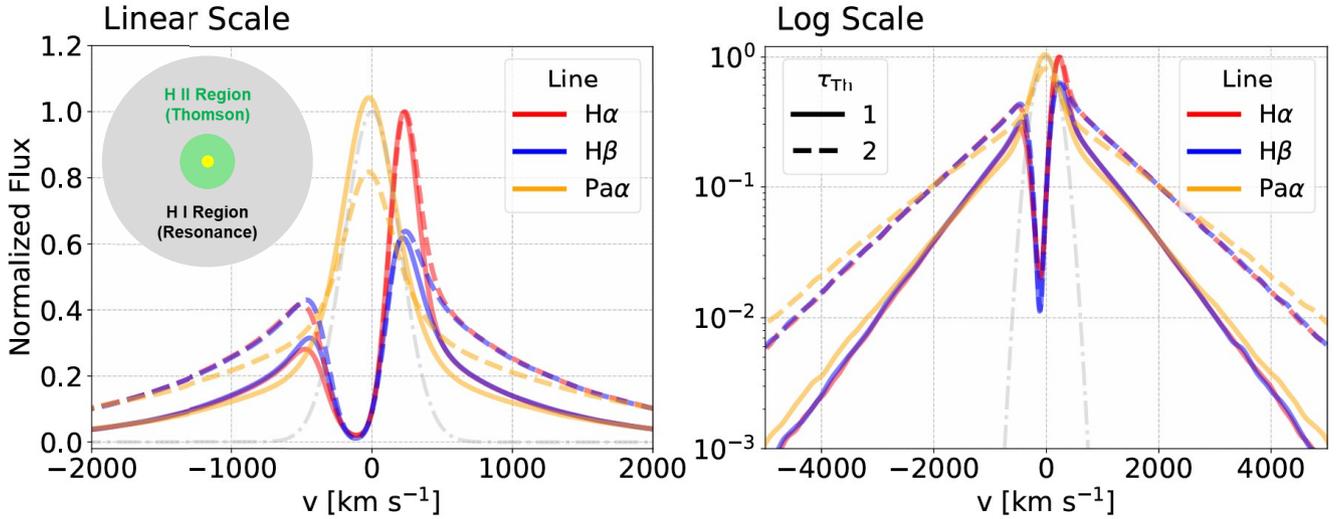


Figure 15. Spectra of $H\alpha$ (red), $H\beta$ (blue), and $Pa\alpha$ (orange) for a model including both Thomson and resonance scattering, shown for two Thomson optical depths: $\tau_{\text{Th}} = 1$ (solid) and 2 (dashed). Fluxes are normalized at $v = 2000 \text{ km s}^{-1}$, assuming a spectral resolution of $R = 3000$ ($\sim 100 \text{ km s}^{-1}$) and an intrinsic Gaussian ($\sigma_{\text{src}} = 200 \text{ km s}^{-1}$; shown in grey). The scattering geometry (the sketch in the left panel) consists of a central source surrounded by a spherical region of radius R , with an inner ionized (H II) region ($r < 0.01R$) producing Thomson scattering, and an outer neutral (H I, with significant $n = 2$ population; $0.01R < r < R$) producing resonance scattering. The H II region has a temperature $T_e = 10^4 \text{ K}$, while the H I region is characterized by $N_{\text{HI}, n=2} = 10^{16} \text{ cm}^{-2}$, an outflow velocity of 100 km s^{-1} , and a random velocity dispersion of 100 km s^{-1} . Left and right panels show linear and logarithmic flux, respectively.

which are high enough to cause significant Raman-scattered features around $H\alpha$ and $H\beta$.

Consequently, without strong emission in the higher Lyman series, Raman scattering cannot be the dominant process responsible for the broad wings around Balmer lines in LRDs. However, if $\text{Ly}\beta$ and $\text{Ly}\gamma$ emission lines are present, Raman scattering could still contribute, producing a characteristic FWHM ratio of $H\alpha$ to $H\beta$ of $\gtrsim 1.28$ (cf. Fig. B2 and Appendix B). Future higher signal-to-noise observations of larger LRD samples may be able to detect such subtle line width differences. Additionally, while higher Lyman-series photons are converted to Balmer lines due to Raman scattering, in practice a fraction of $\text{Ly}\beta$ and $\text{Ly}\gamma$ photons may escape and become observable. Currently, such higher Lyman lines have not been detected in LRDs, but future rest-UV spectroscopy (e.g. R. Tripodi et al. 2025) may uncover them. If such Lyman emission features are observed, comparing their measured spectral widths with those required for Raman broadening could serve as a test of this scenario. As discussed above, Raman scattering can explain $H\alpha$ wings of $\sim 1000 \text{ km s}^{-1}$ only if the corresponding $\text{Ly}\beta$ line has an FWHM of $\sim 150 \text{ km s}^{-1}$. Therefore, detecting a higher $\text{Ly}\beta/H\alpha$ width ratio would argue against Raman scattering from higher Lyman-series emission as the primary origin of the observed broad wings.

4.4 Combining Thomson and Resonance scattering

The spectra of hydrogen emission lines in LRDs exhibit broad components of similar width in both Balmer and Paschen lines (ranging from 1000 to 2000 km s^{-1} ; M. Brazzini et al. 2025; I. Juodžbalis et al. 2025), along with absorption-like features near the line centre, but only for the Balmer lines. As shown in Section 3.1, resonance scattering of $H\alpha$ and $H\beta$ produces P-Cygni-like or symmetric double-peaked profiles in Balmer lines, while Thomson scattering contributes similarly broad wings to both Balmer and Paschen lines (Section 3.3). These effects are consistent with several spectral features observed in LRDs. Therefore, in this section, we

explore how the combination of these two scattering mechanisms shapes the emergent hydrogen line profiles.

We adopt a scattering geometry with a central emission source and a spherical scattering medium of radius R . The central source emits hydrogen lines – $H\alpha$, $H\beta$, and $Pa\alpha$ with intrinsic widths of $\sigma_{\text{src}} = 200 \text{ km s}^{-1}$ and flux ratios under Case B recombination. We assume that the intrinsic width is less than the electron thermal motion, $v_{e,\text{th}} (\approx 500 \text{ km s}^{-1}$ at 10^4 K), as the line broadening due to Thomson scattering becomes negligible when the width exceeds $v_{e,\text{th}}$; we will discuss the dependence on the width of the Thomson broadening in Section 4.6.

The scattering medium is composed of two layers: an inner ionized region responsible for Thomson scattering and an outer neutral region causing resonance scattering. The inner region (within $0.01R$) is fully ionized and characterized by electron temperature T_e and Thomson optical depth τ_{Th} . The outer region is a neutral H I gas with the $2s$ population, characterized by column density $N_{\text{HI}, 2s}$, outflow velocity v_{exp} , and turbulent velocity σ_R . Here, we assume a dense and compact H II region, as a H II region would be highly ionized to be Thomson thick. Thus, Thomson-scattered photons in this compact region act like the broad incident radiation from the central source. Note that the formation of $H\alpha$ and $H\beta$ line profiles depends on the radius of the H II region relevant for Thomson scattering (see Appendix D). In addition, if an outflow velocity of H II region is comparable to or exceeding the electron thermal speed, the Thomson-scattered wings become asymmetric (Appendix C, see also L. H. Auer & D. van Blerkom 1972; C. Huang & R. A. Chevalier 2018).

Intrinsic photons emitted from the central source first undergo Thomson scattering in the inner H II region. Then, $H\alpha$ and $H\beta$ photons experience additional resonance scattering in the outer H I layer. In contrast, $Pa\alpha$ photons do not experience resonance scattering, as the $n = 3$ level has a much shorter lifetime than the $n = 2$ state (see Fig. 2) and hence is unlikely to be overpopulated.

Fig. 15 shows the emergent spectra of $H\alpha$, $H\beta$, and $Pa\alpha$ for $\tau_{\text{Th}} = 1$ and 2, which correspond to electron column density $\approx 1.5 \times 10^{24}$ and $3.0 \times 10^{24} \text{ cm}^{-2}$, respectively. Both $H\alpha$ and $H\beta$ exhibit P-Cygni-like

profiles caused by resonance scattering and develop broad wings due to Thomson scattering. Pa α shows a single-peaked profile centred at zero velocity, consistent with no resonance scattering. In the left panel (linear scale), the core profiles of H α and H β ($|v| < 500 \text{ km s}^{-1}$) do not depend on τ_{Th} . In the right panel (log scale), the line wings ($|v| > 500 \text{ km s}^{-1}$) are significantly enhanced with increasing τ_{Th} , demonstrating the dominant role of Thomson scattering in producing broad wings in this case.

Interestingly, the red peak of the emergent H α is more prominent than that of H β (cf. left panel of Fig. 15). This is because multiple resonance scatterings convert H β photons into Pa α and H α (as discussed Section 3.1.3). This enhanced red peak of H α with respect to H β is also seen in observations (I. Juodžbalis et al. 2024; M. Brazzini et al. 2025), indicating the importance of Balmer resonance scattering.

Another noteworthy spectral feature is the similarity of the absorption features in the blue parts of the spectra (cf. Fig. 15). This similarity is not seen in observations, where typically the H α absorption-like feature is weaker than that of H β . This discrepancy likely results from the fact that H α emission may originate not only from regions embedded in optically thick gas but also from additional regions less affected by strong scattering. The absence of a broad component of [O III] forbidden line (I. Juodžbalis et al. 2024, 2025; X. Ji et al. 2025a, b; R. Maiolino et al. 2025a), suggests that partial hydrogen emission originates from regions not enshrouded by optically thick gas.² If these outer regions produce hydrogen lines via collisional excitation and have a high H α /H β ratio, the resulting H α profile will show weaker absorption. This additional emission can ‘fill’ the line centre, thus producing dissimilar spectral features.

Quantifying the (dis)similarity in the observations, and the comparison to detailed multifrequency RT models will help to characterize the scattering medium of LRDs and help to disentangle the potentially multiple emission sources. In future work, we will expand this study using a larger simulation library to fit observed LRD spectra, considering different scattering geometries and additional emission sources. Furthermore, the high-resolution spectroscopy of Balmer lines in LRDs is required to understand their formation.

4.5 Formation of hydrogen Balmer lines in optically thick, dense gas

Although our simulations only consider resonance scattering of two Balmer lines, combining Thomson and resonance scattering already produces a wide diversity of line profiles. Under the Case B recombination assumption (on-the-spot approximation), higher Lyman-series lines (Ly β , Ly γ , etc.) are assumed to be optically thick and to convert into lower series lines via interaction with neutral hydrogen. If the H I gas has a high $n = 2$ population and is optically thick to all Balmer transitions (i.e. on-the-spot approximation for Balmer lines), only H α photons can escape, due to the metastable nature of the 2s state. This modifies the Balmer line ratios. Therefore, various H α /H β ratios observed in LRDs (M. Brooks et al. 2025; A. J. Taylor et al. 2025b) may be due to RT effects, rather than dust extinction.

The collisional process in a dense gas, such as the BLR in AGN and stellar atmosphere, affects the formation of Balmer lines (S. A. Drake & R. K. Ulrich 1980; S. Collin-Souffrin, S. Dumont &

J. Tully 1982; K. T. Korista & M. R. Goad 2004; G. J. Ferland et al. 2017). At a high electron number density ($n_e > 10^8 \text{ cm}^{-3}$), the collisional de-excitation is not negligible, thus electrons in the $n = 4$ states are easier de-excited than in the $n = 3$ state. It causes a difference between intrinsic H α and H β profiles, and their emission region is not identical. Thus, if the Thomson scattering originates from a dense gas (e.g. M. C. Begelman & J. Dexter 2025), H α and H β emission experience Thomson scattering with different optical depth, resulting in a difference in strength of the broad component around the hydrogen emission lines (M. Brazzini et al. 2025).

This extreme gas condition may be relevant not only for LRDs and AGN, but also for other objects where H α shows strong emission while higher Balmer lines show an absorption-dominated feature. In future studies, we will investigate the formation of Balmer lines under these conditions, including the effects of scattering on higher order Balmer lines such as H γ and H δ .

4.6 Overestimating M_{BH} due to Thomson scattering

Emission-line profiles in AGN spectra are essential for estimating the SMBH mass M_{BH} with the naive expectation being $M_{\text{BH}} \propto \text{FWHM}^2$. Using samples of (low-redshift) black holes where the mass is precisely determined through reverberation mapping, as well as the Balmer line widths are measured, one can calibrate the $M_{\text{BH}}-\text{FWHM}$ relation empirically. This has been done in numerous studies (J. E. Greene & L. C. Ho 2005; A. E. Reines & M. Volonteri 2015; J.-H. Woo et al. 2015). Specifically, A. E. Reines & M. Volonteri (2015) found

$$\log M_{\text{BH}} = 6.60 + 0.47 \log \left(\frac{L_{\text{H}\alpha}}{10^{42} \text{ erg s}^{-1}} \right) + 2.06 \log \left(\frac{\text{FWHM}_{\text{H}\alpha}}{10^3 \text{ km s}^{-1}} \right) \quad (17)$$

where $L_{\text{H}\alpha}$ and $\text{FWHM}_{\text{H}\alpha}$ are the luminosity and FWHM of the H α broad component, respectively.

However, as mentioned above, this relation is calibrated using a local (mostly $z < 0.055$) AGN sample and its applicability to LRDs – presumably AGN at high redshift – is still uncertain. Moreover, Balmer absorption features commonly seen in LRDs are rare in local AGNs (P. B. Hall 2007; X. Lin et al. 2025), suggesting denser or more complex gas environments. Additionally, the exponential shape of the broad wings observed in LRDs (which is typically absent in local AGNs) suggests Thomson scattering may contribute. Therefore, it is important to evaluate how Thomson scattering could bias M_{BH} measurements in such systems.

In this section, we assess how Thomson scattering affects black hole mass estimates under two scenarios: (1) a Thomson-dominated case, and (2) a Thomson-thick BLR case. These two setups are illustrated in Fig. 16. Note that a similar analysis was recently carried out by V. Rusakov et al. (2025) who assumed that a narrow-line emission region is embedded within Thomson-thick gas. Here, we take a more agnostic approach to the intrinsic emission mechanism of the Balmer lines.

4.6.1 Thomson-dominated case

In the first scenario, the Thomson-dominated case, we assume that the broad component of H α originates from only Thomson scattering. The H α emission region is surrounded by free electrons (i.e. an H II region). The intrinsic emission undergoes Thomson scattering and develops broad wings, which are similar to the spherical geometry explored in Section 3.3. We consider two intrinsic line widths,

²Note that the broad wing feature and P-Cygni-like profile of He I line at 10 830 Å observed in LRDs (I. Juodžbalis et al. 2024; X. Lin et al. 2025) can also be due to scattering processes.

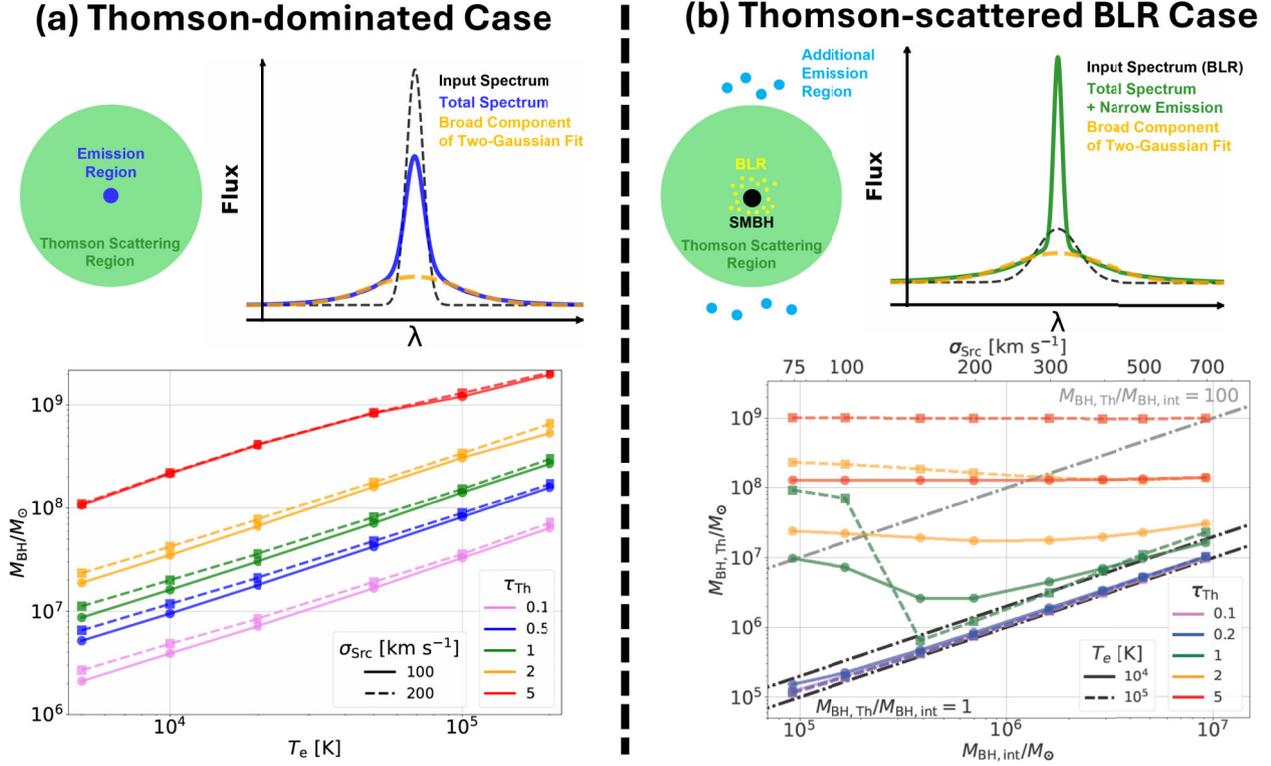


Figure 16. Schematic illustrations (top) and black hole mass estimates (bottom) for two scenarios involving Thomson scattering. Left: *Thomson-dominated case* – a central source embedded in an ionized, optically thick medium produces electron-scattered profiles with exponential wings. The resulting spectra are fit with two Gaussians to estimate M_{BH} as a function of electron temperature and optical depth. Right: *scattered-BLR case* – a virialized BLR emits through a Thomson-scattering in H II region, with an added narrow-line component. The bias in M_{BH} due to scattering is shown as a function of intrinsic mass. See the text for model details.

$\sigma_{\text{src}} = 100$ and 200 km s^{-1} , and run simulations in a range of electron temperatures $T_e = 5 \times 10^3 - 2 \times 10^5 \text{ K}$ and Thomson optical depths $\tau_{\text{Th}} = 0.1 - 5$ (corresponding to electron densities of $n_e \approx 5 \times 10^4 - 2.5 \times 10^6 (R/10 \text{ pc})^{-1}$).³

Although the system contains no actual black hole as the broad component originates from Thomson scattering, not AGN BLR, we fit the emergent line profiles using two Gaussian components to estimate an artificial black hole mass M_{BH} . The top-left panel of Fig. 16 shows an example of this fitting process. The input spectrum (black dashed) is a narrow Gaussian line with width σ_{src} , while the total emergent spectrum (blue solid) includes a broad wing from Thomson scattering. We fit the total spectrum using a two-Gaussian model

$$F_{\text{GG}}(v) = F_{\text{G1}} + F_{\text{G2}} = \frac{A_1}{\sqrt{2\pi}\sigma_1^2} \exp\left(-\frac{v^2}{2\sigma_1^2}\right) + \frac{A_2}{\sqrt{2\pi}\sigma_2^2} \exp\left(-\frac{v^2}{2\sigma_2^2}\right), \quad (18)$$

where F_{G1} and F_{G2} represent the narrow and broad components, respectively, and we impose $\sigma_2 > \sigma_1$ in the fitting process. An example fit for the broad component F_{G2} is shown as the orange dashed line in the top part of Fig. 16.

³As LRDs exhibit a blackbody temperature $\approx 5000 - 6000 \text{ K}$, and the dense gas emitting blackbody radiation is proposed as a Thomson-scattering medium (M. C. Begelman & J. Dexter 2025), we consider a lower electron temperature ($< 10^4 \text{ K}$).

To estimate the black hole mass M_{BH} from the broad spectral component, we fix the total line luminosity $L_{\text{H}\alpha} = 10^{42} \text{ erg s}^{-1}$ consistent with typical LRDs ($L_{\text{H}\alpha} \sim 10^{42-43} \text{ erg s}^{-1}$, J. Matthee et al. 2024; I. Juodžbalis et al. 2025). The luminosity of the broad component is then $A_2/(A_1 + A_2) \times L_{\text{H}\alpha}$. Assuming a Gaussian, we calculate the FWHM of the broad component as $2.355 \sigma_2$. Combining this with equation (4.6), the black hole mass estimate becomes:

$$\log M_{\text{BH}} \approx 6.60 + 0.47 \log\left(\frac{A_2}{A_1 + A_2}\right) + 2.06 \log\left(\frac{2.355 \sigma_2}{10^3 \text{ km s}^{-1}}\right). \quad (19)$$

We apply this equation across simulations with various T_e and τ_{Th} , for both $\sigma_{\text{src}} = 100$ and 200 km s^{-1} .

The bottom-left panel of Fig. 16 shows the resulting inferred M_{BH} values. Note that since there is no intrinsic broad component in this model, there is no black hole and thus no correct black hole mass. The inferred M_{BH} increases with both T_e and τ_{Th} , while dependence on σ_{src} is negligible. When $T_e > 10^5 \text{ K}$ or $\tau_{\text{Th}} > 1$, M_{BH} exceeds $10^7 M_{\odot}$, despite no intrinsic virial motion. This result demonstrates that Thomson scattering alone can mimic a broad component, leading to significant overestimation of M_{BH} through Gaussian fitting. For reference, if the narrow component represents emission from a genuine BLR, an intrinsic velocity of $\sigma_{\text{src}} = 100 \text{ km s}^{-1}$ would correspond to $M_{\text{BH,int}} \approx 10^5 M_{\odot}$. Therefore, if the broad Balmer components of LRDs are shaped by Thomson scattering in ionized gas ($T_e > 10^4 \text{ K}$; V. Rusakov et al. 2025), the inferred M_{BH} may be overestimated by a factor of > 10 .

4.6.2 Thomson-scattered broad-line region case

In the second scenario – the Thomson-scattered BLR case – we consider a geometry similar to the first, with a central emission region surrounded by an H II medium. However, the emission region is treated as the BLR of an AGN, where the gas is virialized by a central SMBH.

The top-right panel of Fig. 16 shows an example spectrum for this setup. To cover a wide range of intrinsic M_{BH} , we adopt $\sigma_{\text{src}} = 75\text{--}700\text{ km s}^{-1}$ (black dashed line), corresponding to $M_{\text{BH,int}} = 10^{4\text{--}7} M_{\odot}$ at fixed $L_{\text{H}\alpha} = 10^{42}\text{ erg s}^{-1}$, based on the M_{BH} scaling relation. We set $T_e = 10^4$ and 10^5 K and explore $\tau_{\text{Th}} = 0.1\text{--}5$. We also include an additional emission region located outside the Thomson-scattering medium, representing a narrow-line region, star-forming region, or outflow. This component has a fixed width of 50 km s^{-1} and contributes half the flux of the intrinsic BLR emission. The total emergent spectrum (green solid line in the top-right panel of Fig. 16) thus consists of Thomson-scattered BLR emission and a narrow Gaussian component. As previously, we fit the total spectrum with two-Gaussian profiles in equation (4.6) and estimate M_{BH} using equation (19), assuming the total H α luminosity fixed at $10^{42}\text{ erg s}^{-1}$.

The bottom-right panel of Fig. 16 shows the resulting M_{BH} estimates $M_{\text{BH,Th}}$ as a function of the intrinsic black hole mass $M_{\text{BH,int}}$. The intrinsic masses are obtained by fitting spectra with no scattering (i.e. $\tau_{\text{Th}} = 0$). The Thomson-scattering fraction is given by $1 - \exp(-\tau_{\text{Th}})$. In the optically thin regime ($\tau_{\text{Th}} < 1$), the scattering fraction is small, and $M_{\text{BH,Th}} \approx M_{\text{BH,int}}$; most scattered photons contribute only weakly to the observed profile. However, even in this case, an exponential wing extending beyond 1000 km s^{-1} is present due to scattering.

In the optically thick regime ($\tau_{\text{Th}} \gtrsim 1$), the scattering fraction becomes significant (≈ 0.86 at $\tau_{\text{Th}} = 2$), and Thomson scattering dominates the broad component formation. This leads to a nearly flat $M_{\text{BH,Th}}$ curve, with mass overestimation by more than an order of magnitude. At intermediate optical depth ($\tau_{\text{Th}} \approx 1$), the overestimation depends on the intrinsic line width. For narrow intrinsic width ($\sigma_{\text{src}} < 200\text{ km s}^{-1}$), the scattered photons are interpreted as the primary broad component. For a broader width ($\sigma_{\text{src}} > 200\text{ km s}^{-1}$), the scattering causes additional broadening, leading to M_{BH} overestimates by factors of ~ 2 . Consequently, if Thomson scattering occurs in the gas surrounding the BLR, the observed line width can be significantly broadened by thermal electron motions, resulting in overestimated black hole masses.

4.6.3 Implications for line profiles and M_{BH} estimates

In summary, this section investigates the effect of Thomson scattering on black hole mass measurements. We explored two scenarios: (1) a Thomson-dominated wing and (2) a Thomson-scattered BLR. In both cases, we estimated M_{BH} using the standard scaling relation calibrated from reverberation mapping of local AGN. In the first case, we showed that pure Thomson scattering can generate broad wings around Balmer lines, resulting in fictitious black hole masses exceeding $10^6 M_{\odot}$ when $T_e \geq 10^4\text{ K}$ and $\tau_{\text{Th}} \geq 0.1$. In the second case, we demonstrated that scattering of BLR emission broadens the observed profile and leads to overestimation of M_{BH} , particularly when $\tau_{\text{Th}} \geq 1$. Although Thomson scattering has a weaker impact on M_{BH} estimates in the optically thin regime ($\tau_{\text{Th}} < 1$), it still produces exponential wings in the far line wings, as seen in local AGNs (e.g. A. Laor 2006).

The role of Thomson scattering in LRDs remains debated. V. Rusakov et al. (2025) proposed that the broad H α features in LRDs

are dominated by Thomson scattering, suggesting that intrinsically narrow emission lines with widths $\sim 100\text{--}200\text{ km s}^{-1}$ are broadened through scattering in dense, ionized gas, which is similar to our first scenario. In contrast, M. Brazzini et al. (2025) argued that the broad hydrogen line features in LRDs are inconsistent with this Thomson-dominated scenario. They suggest that if scattering dominates, the H α -scattering fraction should be higher than that of H β due to dust extinction. However, as discussed in Section 4.5, the Thomson optical depth of H β can be higher than that of H α in dense environments, complicating this interpretation. These contrasting views highlight the need for detailed RT modelling to understand the formation of broad emission features in LRDs.

Our scenarios have important implications for LRDs, where Balmer absorption features suggest dense gas environments. Thomson scattering can naturally produce exponential wings and apparent line broadening, potentially biasing M_{BH} estimates. In addition, if the bulk velocity of the scattering medium is faster than the electron thermal speed, Thomson scattering can provide asymmetric wings on emission lines, depending on the velocity structure of the gas. However, line profile analysis alone is insufficient to distinguish between kinematic and scattering origins of broad wings. Since Thomson scattering induces strong polarization (H.-W. Lee 1999; H. J. Kim et al. 2007), future spectropolarimetric observations of LRDs are a potential avenue to test this hypothesis.

5 CONCLUSION

In this study, we explored the physical origin of the hydrogen line profiles observed in LRDs, which are compact, high-redshift sources identified through *JWST* spectroscopy. These systems exhibit broad wings (FWHM $\gtrsim 1000\text{ km s}^{-1}$) in Balmer and helium emission lines, along with Balmer absorption features and a strong Balmer break, indicating a dense, neutral hydrogen medium in the $n = 2$ state. Under such conditions, RT effects may affect the formation of hydrogen lines.

We performed 3D Monte Carlo RT simulations to examine three key scattering mechanisms that shape hydrogen line profiles: (1) resonance scattering by hydrogen in the $n = 2$ state, (2) Raman scattering of UV photons by ground-state hydrogen ($n = 1$), and (3) Thomson scattering by free electrons. For H β , we explicitly include multibranching transitions in the resonance scattering process, accounting for both $n = 4 \rightarrow 2$ and $n = 4 \rightarrow 3$ decay channels. These pathways allow H β photons to convert into Pa α and H α , significantly suppressing the emergent H β flux. For Raman scattering, we adopt wavelength-dependent cross-sections and branching ratios near the Lyman series. For Thomson scattering, we explore a range of electron temperatures (T_e) and optical depths (τ_{Th}).

Our results from each scattering process show that:

(i) Resonance scattering explains the distinct profile differences between H α , H β , and Pa α and leads to photon conversion from H β to Pa α and H α , altering flux ratios (i.e. leading to large deviations from the value under the Case B recombination, see Fig. 5) and absorption depths. We have demonstrated how these effects imprint the gas geometry and kinematics on the emergent line shapes (Section 3.1), and will explore their diagnostic power systematically in future work.

(ii) While resonant scattering can significantly alter (and for very large optical depths) widen the emergent H α spectrum significantly, this is not the case for H β due to the efficient conversion to Pa α (Section 4.2).

(iii) Raman scattering can produce extremely broad wings (up to $\sim 10^4\text{ km s}^{-1}$), with widths scaling approximately as $\text{FWHM} \propto \sqrt{N_{\text{HI}}}$

and different degrees of broadening for each Balmer line (Fig. 11 and Section 3.2). We show that the observed Balmer line widths are not consistent with these continuum Raman-scattered features. In addition, the UV continuum in LRDs is likely too faint, effectively ruling out Raman-scattered continuum photons to be the origin of the broad Balmer lines observed (Section 4.3).

(iv) However, we also show that if higher Lyman series with the width $\sim 150 \text{ km s}^{-1}$ are fully Raman scattered, the width of the Raman wing around $H\alpha$ and $H\beta$ is over 1000 km s^{-1} , with an FWHM ratio of $H\alpha$ and $H\beta$ is ≈ 1.28 (Fig. B2 and Appendix B).

(v) Thomson scattering produces symmetric broad wings with similar widths across all hydrogen lines. In the optically thin regime, the wings exhibit exponential profiles, with widths scaling as FWHM $\propto \sqrt{T_e}$ (Figs 11 and 12). At higher optical depths ($\tau_{\text{Th}} \gtrsim 2$), the wings deviate from the exponential shape due to multiple scatterings, and their width further increases with τ_{Th} (Figs 12 and 13).

We show that a combination of Thomson and resonant scattering can explain most spectral features observed in LRDs, specifically the broad wings and the central asymmetric double peak (Section 4.4). However, to understand different absorption features and broad wings of Balmer lines, the additional RT effects, such as on-the-spot approximation of Balmer lines, collisional excitation, and collisional de-excitation, are required (Section 4.5).

Importantly, we demonstrate that if the broad component of Balmer lines is shaped by Thomson scattering rather than intrinsic broad emission, then black hole masses estimated using methods for local AGNs can be significantly overestimated (Section 4.6). For optically thick cases ($\tau_{\text{Th}} \geq 2$), the overestimation can exceed a factor of 10 (Fig. 16).

Our results show the importance of detailed RT modelling when interpreting hydrogen line profiles, not only for LRDs and AGN, but also for other astrophysical systems exhibiting broad hydrogen wings and absorption features. In future work, we will apply our simulation library to analyse observed spectra of LRDs and expand our models to include more complex environments, such as those combining resonance and Thomson scattering in multiphase media.

ACKNOWLEDGEMENTS

The authors thank the anonymous referee for constructive comments, which improved the clarity of this paper. SJC acknowledges support from the ERC synergy grant 101166930 – RECAP. MG thanks the Max Planck Society for support through the Max Planck Research Group, and the European Union for support through ERC-2024-STG 101165038 (ReMMU). JM acknowledges funding by the European Union (ERC, AGENTS, 101076224). CAM acknowledges support by the European Union ERC grant RISES (101163035), Carlsberg Foundation (CF22-1322), and VILLUM FONDEN (37459). Computations were performed on HPC systems Freya and Orion at the Max Planck Computing and Data Facility.

DATA AVAILABILITY

The data of this study are available upon reasonable request to the first author.

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APPENDIX A: RAMAN-SCATTERING CROSS-SECTION

Fig. A1 shows the total scattering cross-section for both Rayleigh and Raman scatterings near $\text{Ly}\beta$, $\text{Ly}\gamma$, and $\text{Ly}\delta$, as well as the branching ratios for Raman scattering to $n = 2, 3$, and 4 states, computed using the Kramer–Heisenberg formula (H.-W. Lee 2013; S.-J. Chang et al. 2015).

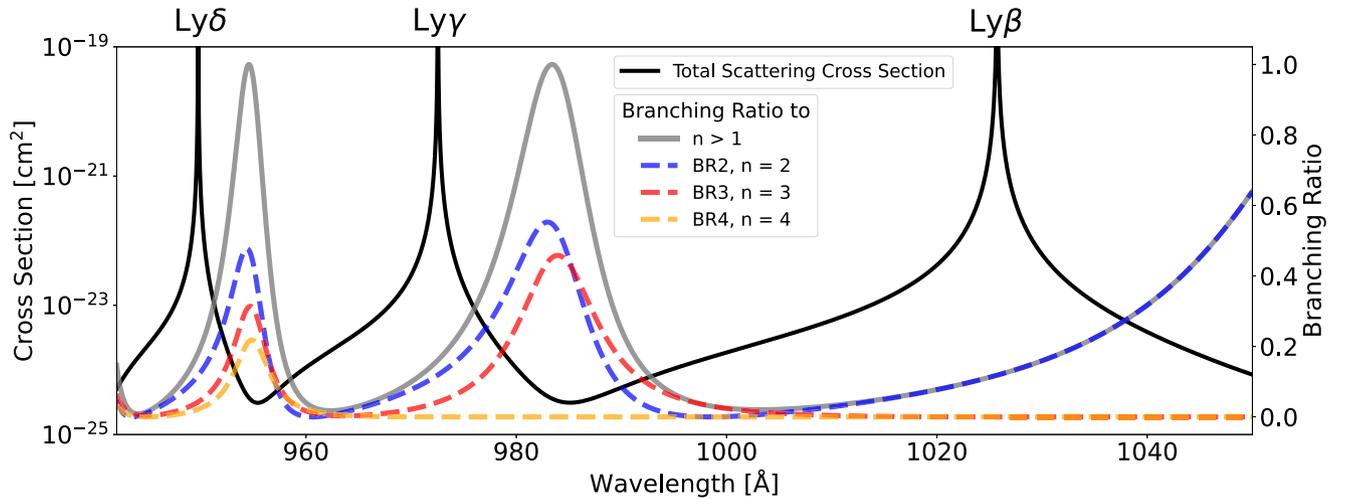


Figure A1. Scattering cross-section and branching ratios of UV radiation near $\text{Ly}\beta$, $\text{Ly}\gamma$, and $\text{Ly}\delta$ as functions of wavelength. The black solid line represents the total scattering cross-section computed using the Kramer–Heisenberg formula. The coloured dashed lines show the branching ratios for Raman scattering to $n = 2$ (blue), $n = 3$ (red), and $n = 4$ (orange) states.

APPENDIX B: RAMAN BROAD FEATURES BY LYMAN-SERIES EMISSION

As discussed in Section 3.2 (cf. Fig. 11), Raman-scattered H α wings are typically ~ 3 times broader than those of H β when the input radiation is a flat UV continuum. This difference arises from the wavelength-dependent scattering cross-sections near Ly β and Ly γ . However, when the input radiation contains strong Ly β and Ly γ emission lines, the width ratio between Raman H α and H β features can be significantly lower. In particular, if both lines are fully converted into Raman-scattered photons at high N_{HI} , the resulting widths are determined by the intrinsic line profile and the Raman wavelength shift, as given by equation (7). In this case, the width ratio between H α and H β converges to ~ 1.28 .

Fig. B1 shows Raman-scattered H α and H β from Lyman-series emissions for two representative column densities: 10^{21} and 10^{23} cm $^{-2}$. We assume a Gaussian input profile for both Ly β and Ly γ with $\sigma_{\text{src}} = 200$ km s $^{-1}$. At low N_{HI} , the H α Raman wing is $\sim 2 - 3$ times broader than that of H β due to cross-section differences,

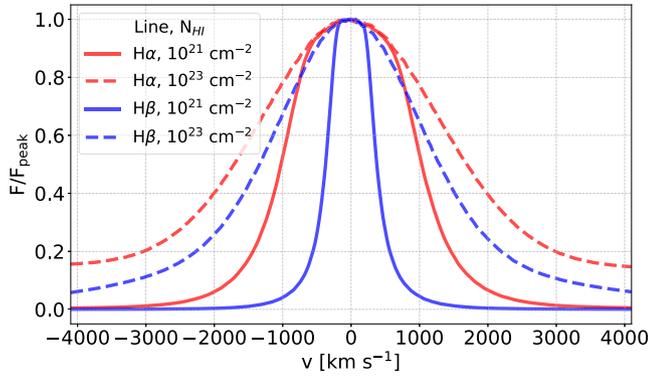


Figure B1. Raman-scattered features around H α (red) and H β (blue) produced by Ly β and Ly γ emission lines. The input Lyman-series lines are assumed to have intrinsic widths of 200 km s $^{-1}$ and an equivalent width of 10 Å. Solid and dashed lines correspond to low (10^{21} cm $^{-2}$) and high (10^{23} cm $^{-2}$) column densities, respectively.

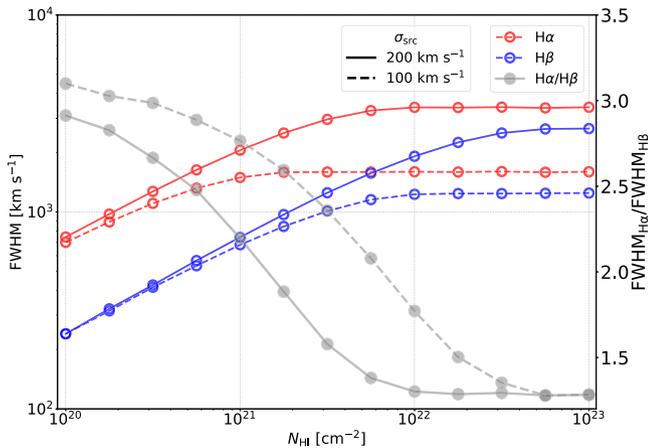


Figure B2. FWHM of Raman features around H α and H β , and their ratio, as a function of N_{HI} . At high column density ($\sim 10^{23}$ cm $^{-2}$), where Lyman-series photons are fully converted into Raman features, the H α /H β width ratio approaches the theoretical value of ~ 1.28 (from equation 7). At low N_{HI} ($\sim 10^{20}$ cm $^{-2}$), the ratio is ~ 3 , reflecting the wavelength dependence of the Raman-scattering cross-section, as seen in the flat continuum case (Fig. 11).

Table B1. Raman broadening factor, $\gamma = \lambda_{\text{sc}}/\lambda_{\text{inc}}$, for various transitions.

Input photon	$N_{\text{HI},100}^*$	Scattered photon	γ
Ly β	$\approx 2 \times 10^{21}$ cm $^{-2}$	H α	6.4
Ly γ	$\approx 10^{22}$ cm $^{-2}$	H β	5.0
		Pa α	19.3
Ly δ	$\approx 5 \times 10^{22}$ cm $^{-2}$	H γ	4.6
		Pa β	13.5
		Br α	42.7

Notes. $*N_{\text{HI}}$ for fully Raman-scattered case. The first and third columns list the incident (Lyman-series) and corresponding Raman-scattered photons, respectively. The second column gives the HI column density $N_{\text{HI},100}$ required for an incident Lyman line with a width of 100 km s $^{-1}$ to be fully Raman scattered. The fourth column shows the velocity broadening factor γ , which relates the FWHM of the scattered and incident profiles.

resembling the flat continuum case shown in Fig. 10. At high N_{HI} , the H α and H β profiles become comparable in width, as nearly all input photons are scattered.

To quantify the trend of Raman H α and H β widths, Fig. B2 shows the FWHM of the Raman H α and H β features, and their ratio, as a function of N_{HI} , for two input widths: $\sigma_{\text{src}} = 100$ and 200 km s $^{-1}$. At low N_{HI} ($< 10^{21}$ cm $^{-2}$), the scattering cross-section primarily controls the profile width and favours broader H α wings. As N_{HI} increases, the conversion probability becomes saturated, and the intrinsic width σ_{src} begins to dominate the emergent profile. At high N_{HI} ($> 10^{22}$ cm $^{-2}$), the H α /H β FWHM ratio converges to the theoretical value of ~ 1.28 , determined by the Raman wavelength shifts.

When Lyman-series emission lines of equal intrinsic width are fully Raman scattered, the width of the resulting Raman features is set by the broadening factor in equation (7). As noted in Section 2.3.1, the additional broadening in velocity space is

$$\gamma = \frac{\Delta v_{\text{sc}}}{\Delta v_{\text{inc}}} = \frac{\lambda_{\text{sc}}}{\lambda_{\text{inc}}}, \quad (\text{B1})$$

where λ_{sc} and λ_{inc} are the wavelengths of the scattered and incident photons, respectively. Table B1 lists γ for each transition, along with the HI column density $N_{\text{HI},100}$ at which an incident Lyman line of 100 km s $^{-1}$ width is fully Raman scattered. For column densities below $N_{\text{HI},100}$, the FWHM ratio between Raman H α and H β reverts to the UV continuum case (≈ 3 as shown in Fig. 11). Notably, since both Raman H β and Pa α originate from an inelastic scattering of Ly γ , their FWHM ratio is fixed at ≈ 3.86 , which is the ratio of their respective broadening factors.

In summary, Raman scattering of Ly β and Ly γ emission lines leads to H α and H β broad wings with similar widths when the HI column density is high, but different widths of Balmer and Paschen lines. This effect can mimic the symmetric wing profiles produced by Thomson scattering and must be considered when interpreting the origin of broad components in Balmer lines.

APPENDIX C: THOMSON SCATTERING IN OUTFLOWING GAS

As discussed in Section 3.3, Thomson scattering in a static medium produces symmetric broad wings around emission lines, with a width proportional to $\sqrt{T_e}$. However, when the scattering medium exhibits a bulk outflow comparable to or exceeding the electron thermal velocity, $v_{\text{th,e}} \approx 548$ km s $^{-1}$ at $T_e = 10^4$ K, Doppler shifts by the bulk motion can break the symmetry of the Thomson-scattered wings (e.g. L. H. Auer & D. van Blerkom 1972; C. Huang & R. A. Chevalier 2018). For this reason, we consider an outflow in

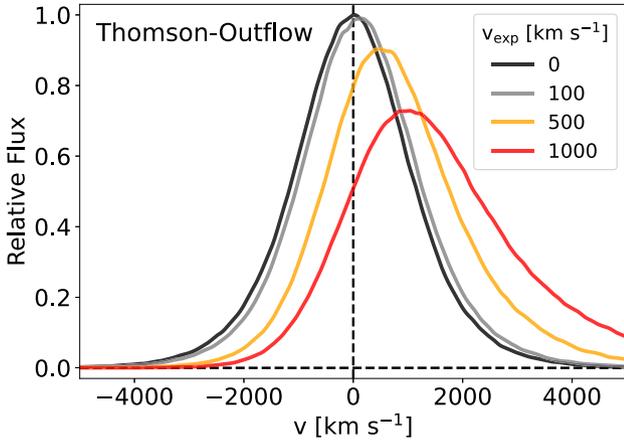


Figure C1. Spectra composed of Thomson-scattered photons for various outflow velocities v_{exp} , normalized by the peak flux of the static case ($v_{\text{exp}} = 0 \text{ km s}^{-1}$). The line colours correspond to $v_{\text{exp}} = 0$ (black), 100 (grey), 500 (orange), and 1000 km s^{-1} (red). As v_{exp} increases, the spectra become asymmetric, with enhanced red wings.

the spherical H II geometry for Thomson scattering and assume a constant radial outflow velocity, which is similar to the outflow velocity in the geometry for resonance scattering in Section 2.2.

Fig. C1 shows Thomson-scattered spectra for various outflow velocities, $v_{\text{exp}} = 0\text{--}1000 \text{ km s}^{-1}$, at an electron temperature $T_e = 10^4 \text{ K}$ and Thomson optical depth $\tau_{\text{Th}} = 1$. At $v_{\text{exp}} \lesssim 100 \text{ km s}^{-1}$, the wings remain nearly symmetric since the bulk velocity is much smaller than $v_{\text{th,e}}$. At higher velocities ($v_{\text{exp}} \gtrsim 500 \text{ km s}^{-1}$), the emergent spectrum develops a pronounced red asymmetry, reflecting the Doppler shift of photons scattered by outflowing electrons. This effect becomes significant as $v_{\text{exp}}/v_{\text{th,e}}$ approaches unity or higher, and may serve as a diagnostic of fast, outflowing H II regions in systems where the Balmer wings show measurable red–blue asymmetry. Conversely, when the bulk velocity of the ionized gas is much smaller than $v_{\text{th,e}}$, the scattered wings remain nearly symmetric, as shown in the spectroscopic study of P Cygni (e.g. A. P. Bernat & D. L. Lambert 1978).

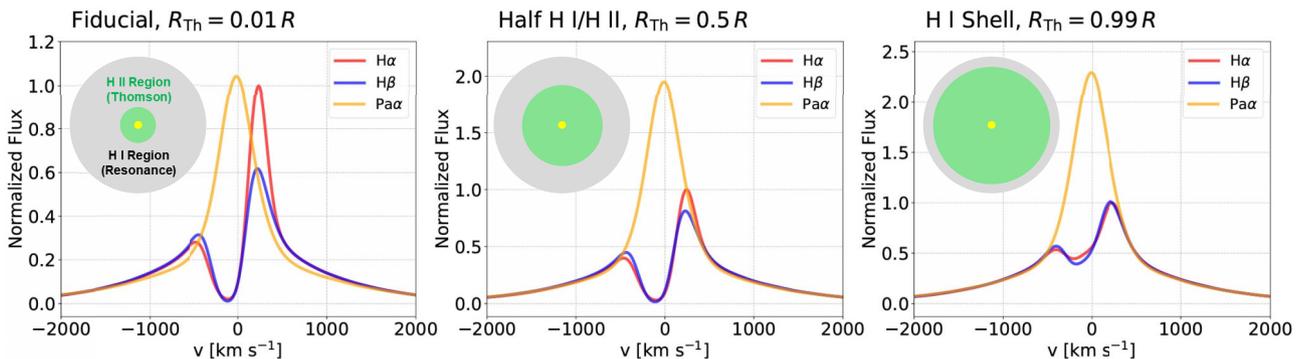


Figure D1. Emergent H α (red), H β (blue), and Pa α (orange) line profiles for different sizes of the Thomson-scattering region. All models assume a spherical geometry with an inner ionized (H II) region for Thomson-scattering and an outer neutral (H I) region with $n = 2$ hydrogen enabling resonance scattering. The flux normalization and other simulation parameters are identical to those used in Fig. 15; all spectra are normalized at $v = 2000 \text{ km s}^{-1}$. From left to right, the Thomson region extends to $R_{\text{Th}} = 0.01R$ (fiducial, matching the $\tau_{\text{Th}} = 1$ model in Fig. 15), $0.5R$, and $0.99R$. As the Thomson-scattering region becomes larger, the broad wings become more prominent, and the differences between the H α and H β profiles driven by resonance scattering are reduced.

APPENDIX D: DEPENDENCE ON RADII OF H II AND H I REGIONS, COMBINING THOMSON AND RESONANCE SCATTERING

In this section, we examine how the size of the inner ionized region for Thomson scattering affects emergent line profiles when both Thomson and resonance scattering are active, as discussed in Section 4.4. Fig. D1 shows the resulting H α , H β , and Pa α spectra for three different Thomson region radii: $R_{\text{Th}} = 0.01R$, $0.5R$, and $0.99R$, all at a fixed Thomson optical depth $\tau_{\text{Th}} = 1$. The outer neutral region, enabling resonance scattering, extends from R_{Th} to R and contains hydrogen in the $n = 2$ state. Other parameters are fixed: outflow velocity $v_{\text{exp}} = 100 \text{ km s}^{-1}$, turbulent velocity $\sigma_{\text{R}} = 100 \text{ km s}^{-1}$, intrinsic width $\sigma_{\text{src}} = 200 \text{ km s}^{-1}$, and H I column density in the $2s$ state $N_{\text{HI},2s} = 10^{16} \text{ cm}^{-2}$, which are identical to those of spectra in Fig. 15.

In Fig. D1, the profile difference between H α and H β diminishes as the Thomson region expands. This trend arises from the interaction between geometry and scattering behaviour. Since the Thomson optical depth is fixed, the emergent energy distribution from the inner H I region remains similar regardless of its size. However, the spatial path taken by photons after emission differs significantly.

In the small Thomson region case (e.g. $R_{\text{Th}} = 0.01R$), photons quickly exit the ionized core and enter the neutral H I shell, where they undergo resonance scattering. Because the H II region is small, photons that scatter in the H I shell are unlikely to re-enter the Thomson region; the effects of resonance and Thomson scattering remain spatially separated. In contrast, when the ionized region is large (e.g. $R_{\text{Th}} = 0.99R$), photons undergoing resonance scattering near the boundary between H I and H II regions can easily scatter back into the ionized zone. These photons are then subject to additional Thomson scattering before escaping. This back-and-forth between the two regions effectively reduces the impact of resonance scattering, as many photons escape via Thomson scattering rather than multiple resonance scattering H I region.

Since the line profile differences between H α and H β originate primarily from resonance scattering, especially due to branching in H β , this effect becomes suppressed as R_{Th} increases. As a result, in the large ionized region case (right panel of Fig. D1), the H α and H β profiles appear nearly identical.

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