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Detection of CH₃⁺ in the O-rich planetary nebula NGC 6302

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ABSTRACT

Planetary nebulae are sites where ejected stellar material evolves into complex molecules, but the precise physical conditions and chemical routes that govern these processes are unclear. The presence of abundant carbon-rich molecules in O-rich environments poses particular challenges. Here we report the first detection of methyl cation (CH₃⁺) in any planetary nebula, observed in the O-rich nebula NGC 6302 using JWST MIRI/MRS

observations. CH_3^+ is a key driver of organic chemistry in UV-irradiated environments. Spatially resolved observations reveal that CH_3^+ is co-located with ^{12}CO , H_2 , H I , HCO^+ , and Polycyclic aromatic hydrocarbons (PAHs). LTE modelling of the CH_3^+ emission yields excitation temperatures of 500-800K in the inner bubble and torus, rising to 1000-2000K in the outer bubble of NGC 6302, with column densities ranging from $\sim 10^{11}$ to 10^{13} cm^{-2} . This detection demonstrates that hydrocarbon radical chemistry must be incorporated into planetary nebulae chemical models. Further near-IR observations are crucial to map different chemical networks operating in these environments.

Keywords: Astrochemistry (75), Circumstellar Gas (238), Planetary nebulae (1249), Small molecules (2267)

1. INTRODUCTION

As of August 2025, more than 330 molecular species have been identified in interstellar and circumstellar environments. A substantial subset of this inventory originates in the circumstellar envelopes of evolved stars, including stars on the asymptotic giant branch (AGB), post-AGB stars, and planetary nebulae (singular: PN; plural: PNe; B. A. McGuire 2022; M. Araki 2025). Elucidating how these molecules form and evolve as local physical conditions change is essential for constraining a wide range of astrochemical and astrophysical processes such as star- and planet formation and the chemical enrichment of galaxies, among others. (see A. G. G. M. Tielens 2013, for a review).

During the AGB phase, low- to intermediate-mass ($\sim 1\text{--}8 M_\odot$) stars undergo severe mass-loss with mass-loss rates ranging from 10^{-8} to $10^{-4} M_\odot \text{ yr}^{-1}$ (S. Höfner & H. Olofsson 2018). These AGB outflows result in the formation of a circumstellar envelope (CSE). As the stellar gas expands radially away from the stellar atmosphere, it cools down, resulting in the formation of molecules and dust. The infrared (IR) spectra of AGB stars indeed show the presence of a rich variety of molecules. More than 90 molecules have been detected in the CSE around carbon-rich stars, and about 27 around oxygen-rich stars (see M. Agúndez et al. 2014; L. Decin et al. 2018; M. Agúndez 2022; B. A. McGuire 2022, and references there in). Note that the molecular species detected in the CSE of AGB stars are mainly simple molecular species, such as CO, C_2H_2 , HCN, CO_2 , H_2O , etc. Eventually, the star progresses through the post-AGB phase and the effective temperature of the star, T_{eff} , increases rapidly with decreasing stellar radii. As $T_{\text{eff}} \gtrsim 20,000\text{K}$, UV radiation from the star ionizes the hydrogen in the previously ejected material and a PN is formed (K. B. Kwitter & R. B. C. Henry 2022). UV photons from the star can also lead to dissociation of molecules in the CSE, and not all molecules survive the transition from the AGB to the PN phase. In fact, a sharp drop is seen in the molecular inventory of PN compared to that of the AGB phase (M. Agúndez 2022). Interestingly, spectral features of larger and more complex molecules like Polycyclic Aromatic Hydrocarbons (PAHs) and fullerenes, such as C_{60} , C_{70} , begin to appear in the post-AGB and PN

phase (M. Matsuura et al. 2005, 2014; N. L. J. Cox et al. 2016; J. Cami et al. 2010; Y. Zhang & S. Kwok 2011).

A key concept in the chemistry of evolved stars is that the chemistry in the AGB outflows depends primarily on the C/O ratio of the stellar photosphere. All stars start their lives as oxygen-rich ($\text{C/O} < 1$) stars, but third dredge-up events in low- and intermediate-mass stars ($\sim 1.5\text{--}4 M_\odot$; D. A. García-Hernández et al. 2013) elevate this C/O ratio above unity late in the AGB phase, converting the envelope to carbon-rich. After H_2 , the next molecule to form in the stellar photosphere is carbon monoxide (CO), which is a highly stable molecule. CO locks up all the free carbon or oxygen atoms – depending on which species is most abundant – leaving the remaining carbon or oxygen to drive further chemistry. Oxygen-rich stellar envelopes thus exhibit the presence of oxides (such as H_2O , SO, AlOH, etc.), and crystalline silicates (R. J. Sylvester et al. 1999; J. Cami 2002; T. Posch et al. 2002; H. Kim et al. 2010; L. Decin et al. 2010; E. De Beck et al. 2013; L. Decin et al. 2018) while carbon-rich stellar envelopes are characterized by the presence of hydrocarbons, PAHs, fullerenes and amorphous dust (T. I. Hasegawa & S. Kwok 2001; J. Cami et al. 2010; M. Agúndez et al. 2014).

However, A. A. Zijlstra et al. (1991) presented the first evidence of mixed or dual chemistry in a PN - where the spectrum of a young PN (IRAS 07027-7934) showed the simultaneous presence of strong PAH emission as well as 1.6 GHz OH maser line emission. Using observations by the Infrared Space Observatory, more cases of dual chemistry in PNe were reported by L. B. F. M. Waters et al. (1998a,b) and M. Cohen et al. (2002). Furthermore, using Spitzer observations, J. V. Perea-Calderón et al. (2009) showed that out of 26 PNe observed towards the galactic bulge, 21 showed signatures of dual chemistry. A larger survey by L. Stanghellini et al. (2012) showed that 46% of the bulge PNe and 24% of the galactic PNe exhibit dual chemistry. It is not clear whether this dual-chemistry phenomenon arises on the AGB or emerges later as a result of processing during the PN phase. Consequently, our knowledge and understanding of the chemical pathways in late stages of stellar evolution, from the AGB phase to PN, is incomplete. It is important to explore how the inventory of molecules is affected by the dual

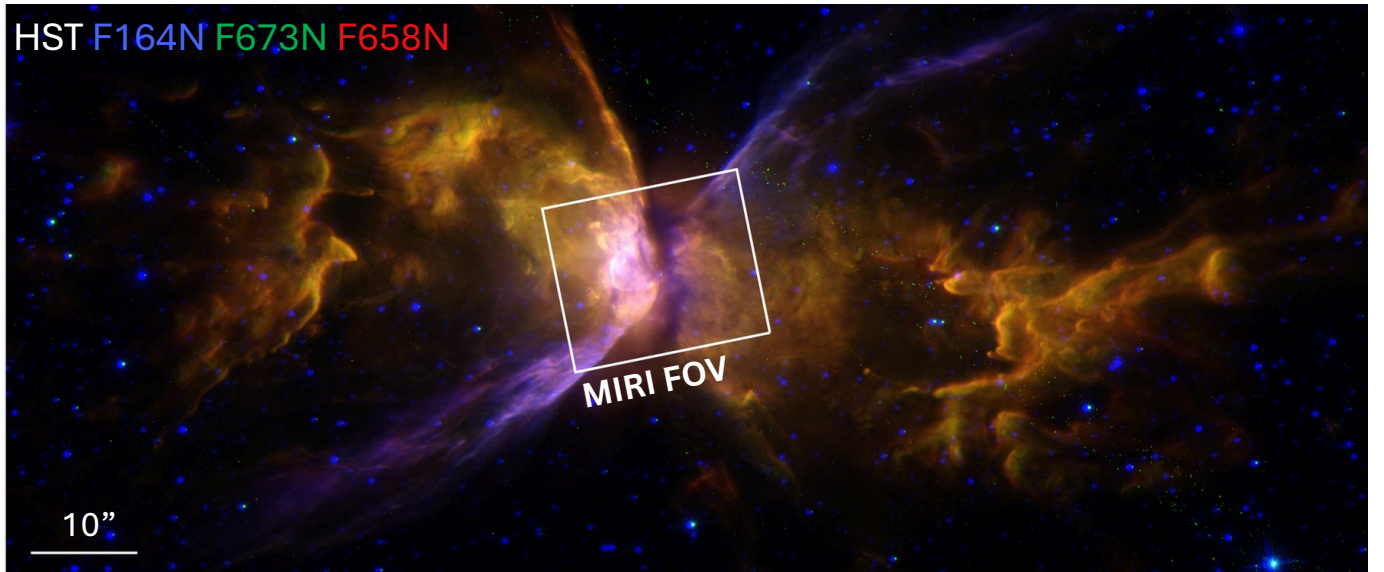


Figure 1. Three-colour composite HST image of NGC 6302, created by Matsuura et al. (2025, MNRAS, submitted) utilizing images from J. H. Kastner et al. (2022) and B. Balick et al. (2023). This image presents a colour overlay of WFC3 observations of NGC 6302, featuring F164N filter in blue, which traces [Fe II] emission, F656N filter in green, which traces H α emission, and F658N in red, which traces [N II] emission. The white box at the center represents the MIRI field-of-view for observations described in Sect. 2.

chemistry, and eventually, what chemical reactions can occur which are not accounted for by strictly C-rich or O-rich chemical pathways. This can potentially provide clues to the formation of other chemical species; for example, the formation of large carbonaceous aromatic molecules like PAHs in an O-rich environment.

In this paper, we report the unambiguous detection of the methyl cation (CH_3^+) in the NGC 6302, a small but important carbonaceous species in what is very clearly an O-rich environment ($\text{C/O} \approx 0.5$, N. J. Wright et al. 2011). While CH_3^+ has long been considered a key species for carbon chemistry in interstellar environments (J. H. Black & A. Dalgarno 1977; D. Smith 1992; A. Sternberg & A. Dalgarno 1995; E. Herbst 2021), its first detection outside the Solar System is a much more recent milestone. CH_3^+ was first detected via JWST/MIRI observations of the irradiated proplyd d203-506 in the Orion Bar by O. Berné et al. (2023). The species is also present at much larger scales in the Orion Bar PDR (M. Zannese et al. 2025) and in the disk of a T-Tauri star (T. Henning et al. 2024). Our detection in NGC 6302 is the first time this species has been detected in a planetary nebula.

2. OBSERVATIONS AND DATA REDUCTION

We use the JWST MIRI/MRS observations of NGC 6302, which were obtained as part of the JWST GO Cycle 1 program 1742. The pointing was centred on the central region of NGC 6302 at R.A. = $17^{\text{h}}:13^{\text{m}}:44.3938^{\text{s}}$, Dec. = $-37^{\circ}:06':12.36$ (J2000). The observations were taken with the MIRI Medium Resolution Spectrometer (MRS) with Integral Field Unit (IFU). The data was obtained for the entire MIRI wavelength range, which is 4.9 to $27.9\mu\text{m}$ spread across 4

channels (M. Wells et al. 2015; I. Argyriou et al. 2023). A mosaic of 5×5 tiles was used to cover the central part of NGC 6302, which includes the central star, the torus, and the bright ionized line emitting region at the innermost region of the bipolar outflow (see fig. 1). Four-point dithering was used to optimize spatial and spectral sampling throughout the field of view.

We used the development version of v1.14.0 JWST Calibration Pipeline with versions 11.17.16 and “jwst.1202.pmap” of the Calibration Reference Data System (CRDS) and CRDS context, respectively. All level 1b files were processed through Detector1Pipeline to obtain level 2a (rate) images. The background exposures were subtracted from the science exposures to get background-subtracted level 2a files. These were then processed through Spec2Pipeline (with residual fringe step tuned on), followed by Spec3Pipeline to produce spectral cube mosaics of NGC 6302. For details on the observations and data reduction, we refer to Matsuura et al., 2025, MNRAS, in press.

3. THE PLANETARY NEBULA NGC 6302

NGC 6302 is a butterfly-shaped planetary nebula, 1.03 ± 0.27 kpc away from Earth (S. Gómez-Gordillo et al. 2020), towards the constellation Scorpius. N. J. Wright et al. (2011) constructed a 3D photoionization model of NGC 6302 and constrained the properties of its central star. They derived an effective temperature of 220,000K, which explains the presence of highly ionized species such as [Mg VII] and [Si VII]. The same study yields a stellar mass of $0.73\text{--}0.82 M_{\odot}$ (with an initial mass of about $5.5 M_{\odot}$), and a luminosity

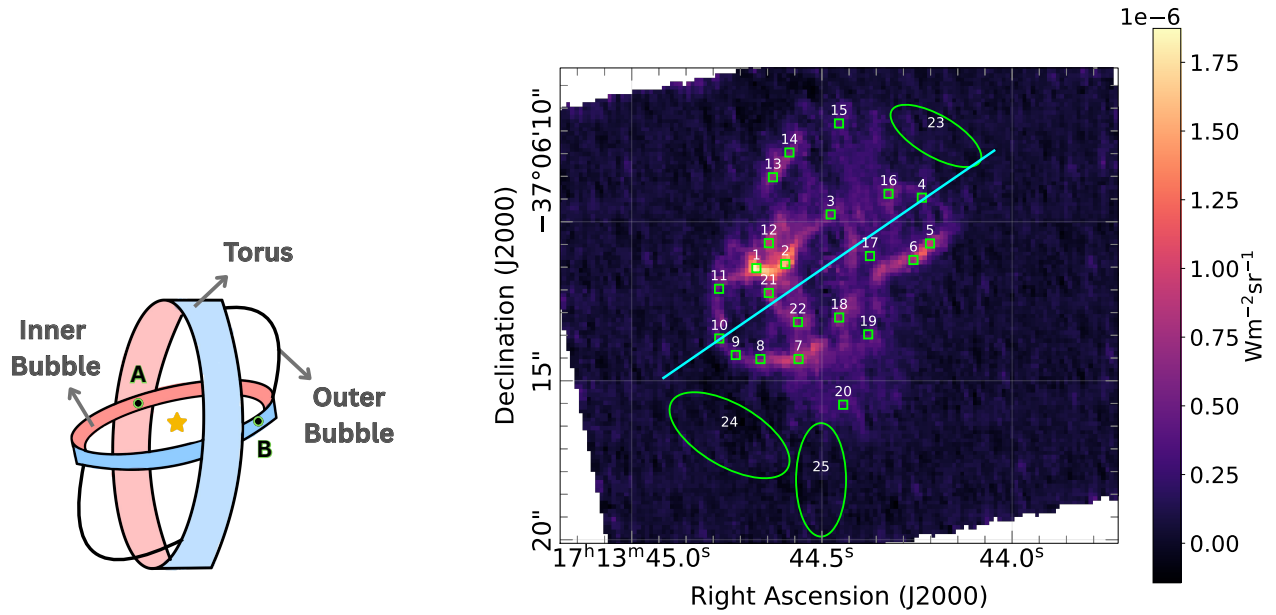


Figure 2. *Left panel:* The cartoon provides a schematic representation of the central region of NGC 6302. Blue and red colours indicate the blueshift and redshift caused by the expansion of the torus and inner bubble from the Earth’s perspective. The schematic is based on the results of [M. Santander-García et al. \(2017\)](#) and [Matsuura et al., 2025, MNRAS, in press](#). The line-of-sight motion of the outer bubble is not constrained and thus is represented by arcs. Points A and B are two bright spots on CH_3^+ integrated surface brightness map. *Right panel:* Integrated surface brightness map of CH_3^+ , obtained by integrating the surface brightness over the $7.13\text{--}7.2\mu\text{m}$ wavelength range at each pixel. The green boxes and ellipses represent the 25 apertures where we analyzed the CH_3^+ emission in more detail. Aperture 1-22 are 2×2 pixels, which is $0.26 \times 0.26''$. Apertures 23-25 are larger elliptical regions with areas of $3.22''^2$, $6.52''^2$, and $4.36''^2$, respectively. The cyan line represents the cut used for making Fig. 9.

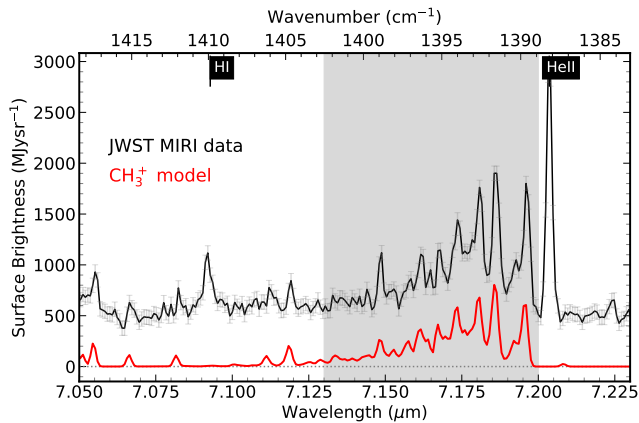


Figure 3. The continuum-subtracted JWST/MIRI spectrum of NGC 6302 at aperture 1 shown in Fig 2 (black, offset vertically for clarity) compared to a CH_3^+ model (red) by [P. B. Changala et al. \(2023\)](#) at excitation temperature of 600 K. The H I recombination line at $7.0928\mu\text{m}$ and the He II line at $7.2036\mu\text{m}$ are labelled at the top.

of $14,300 L_\odot$. NGC 6302 is O-rich with $\text{C/O} \approx 0.5$ ([N. J. Wright et al. 2011](#)).

Optical images (see Fig. 1) reveal two lobes of luminous gas, shaped like the wings of a butterfly. Its bipolar lobes (which stretch as far as 2.1 pc) are exceptionally bright

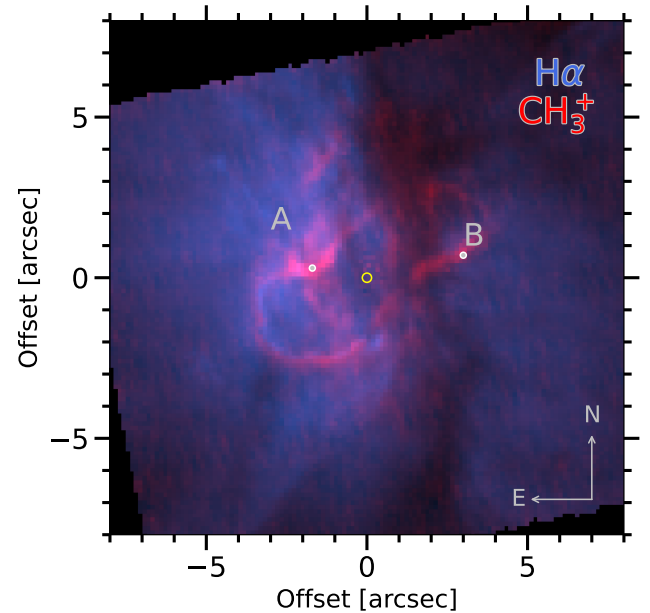


Figure 4. The image represents a colour overlay of HST/WFC3 observations featuring filter F656N in blue ([J. H. Kastner et al. 2022](#)) which traces emission from neutral hydrogen ($\text{H}\alpha$, H I 3–2), and CH_3^+ map in red.

and exhibit a rich variety of emission lines, from the far-ultraviolet to the mid-infrared (S. Casassus et al. 2000; W. A. Feibelman 2001). NGC 6302 is known for exhibiting one of the most complex morphologies seen in PNe. Optical observations (HST/WFC3) also reveal that the two lobes, which are oriented east-west, show small-scale structures such as clumps, tails, knots, and filaments, which seem to be in azimuthally organized zones. [Fe II] emission reveals jets coming out from the central region in south-east and north-west directions, forming an S-shape (J. H. Kastner et al. 2022; B. Balick et al. 2023).

The central region of NGC 6302 comprises three major morphological structures: the *torus* (sometimes referred to as the “central dark lane” or “main ring” in previous studies), the *inner bubble* (sometimes referred to as the “inner ring” in previous studies) and the *outer bubble*. A schematic presentation of these morphological structures is provided in Fig. 2.

The torus is massive and dense with a total (dust and gas) mass of 0.8–3 $M_{\odot}$ and size of 5700 AU (Matsuura et al., 2025, MNRAS, in press). It is the primary reservoir of dust in NGC 6302. These authors point out that the torus is distorted or warped at the edges and is filled with crystalline silicates and other dust grains. In optical observations, the torus appears as a central dark lane that bisects the two lobes and runs nearly north-south, orthogonal to the lobes (see Fig. 1). That dark lane is the front part of the torus that surrounds the central star, obscuring the star in the visible and near-IR (J. H. Kastner et al. 2022). ALMA ^{12}CO J=3–2 maps reveal the torus in great detail and provide kinematic information. The torus is non-Keplerian and is radially expanding at a velocity of 8 km s $^{-1}$. The kinematical age of the torus is estimated to be ~5000–7500 years (N. Peretto et al. 2007; M. Santander-García et al. 2017).

The inner bubble has been observed at optical, mid-IR and sub-mm wavelengths as it is traced by emission of several species such as H I, H $_2$, ^{12}CO , HCO $^+$, PAHs, etc (M. Santander-García et al. (2017); J. H. Kastner et al. (2022), Matsuura et al., 2025, MNRAS, in press, Moraga Baez et al., 2025, under review, Clark et al., in prep). The inner bubble, which runs east-west, is peanut-shaped and is pinched where it intersects with the previously ejected torus (Matsuura et al., 2025, MNRAS, in press). The inner bubble is inclined approximately 60° relative to the torus, and is expanding with a velocity of 11 km s $^{-1}$. It is much younger than the torus, with a kinematical age of ~2200 years (M. Santander-García et al. 2017). B. Balick et al. (2023) describe the sequence of several episodes of ejection from the center of the NGC 6302 in detail.

While the torus and the inner bubble have been detected in optical and sub-mm observations before, JWST MIRI observations by Matsuura et al., 2025, MNRAS, in press revealed two arcs outside the inner bubble, towards the southeast and

northwest, inclined with respect to both the inner bubble and the torus (see Fig. 2). Those arcs constitute the outer bubble, which is bright in H $_2$ and PAH emission. (Matsuura et al., 2025, MNRAS, in press, Clark et al., in prep)

4. CH $_3^+$ EMISSION IN NGC 6302

The JWST/MIRI observations of the central region of NGC 6302 reveal the characteristic CH $_3^+$ emission band in the 7.13–7.20 μm range (see Fig. 3) and show that the emission is spatially extended and resolved across distinct regions.

4.1. The spatial distribution of CH $_3^+$

To reveal the spatial distribution of CH $_3^+$, we integrated the surface brightness over the wavelength range 7.13–7.20 μm after continuum subtraction for each spatial pixel in the MIRI field of view. We restricted the integration to these wavelengths to avoid contributions from other species. An example of the adopted continuum is shown in Appendix A, and the resulting map of the integrated surface brightness is shown in Fig. 2. The map shows bright CH $_3^+$ emission along the inner bubble’s edge (east-west orientation) and the torus (north-south orientation). Although not visible on the map, we found that CH $_3^+$ is also present in the outer bubble (see below), but the column densities are too low to be visible in the integrated surface brightness map. To guide the reader and facilitate spatial comparison, we compared the CH $_3^+$ map to the HST/WFC3 image (through the F656N filter; see J. H. Kastner et al. 2022), which traces H α emission, in Fig. 4. Note that the CH $_3^+$ emission along the torus overlaps with what appears as a central dark lane in the H α observations. Interestingly, no clear signature of CH $_3^+$ is detected along the outer bubble.

To better understand what drives the CH $_3^+$ emission, we investigate its spectroscopic characteristics and spatial variations using defined apertures rather than individual pixels to achieve better signal-to-noise (S/N) ratios. We selected 25 apertures to sample all distinct regions with CH $_3^+$ emission along the torus, inner bubble, and outer bubble (see Fig. 2). Apertures 1–22 use 2 $\times$ 2 pixel areas, each covering 0.26 $\times$ 0.26'' 2 given the 0.13'' pixel scale in channel 1c. Apertures 23–25 are larger elliptical regions (with areas 3.22'' 2 , 6.52'' 2 , and 4.36'' 2 respectively) to achieve adequate S/N for the weaker CH $_3^+$ emission in the outer bubble. These apertures were positioned to cover the outer bubble arc structure as guided by the H $_2$ map (see Appendix B). The aperture distribution is as follows: apertures 1–2 cover the intersection of the inner bubble and torus; apertures 3–6 and 7–11 trace the west and east edges of the inner bubble, respectively; aperture 12 lies north of the intersection point; apertures 13–15 trace the north rear side of the torus; apertures 16–20 and 21–22 trace the front and rear edges of the torus, respectively; and apertures 23–25 sample the outer bubble’s northwest arc and south-southwest region.

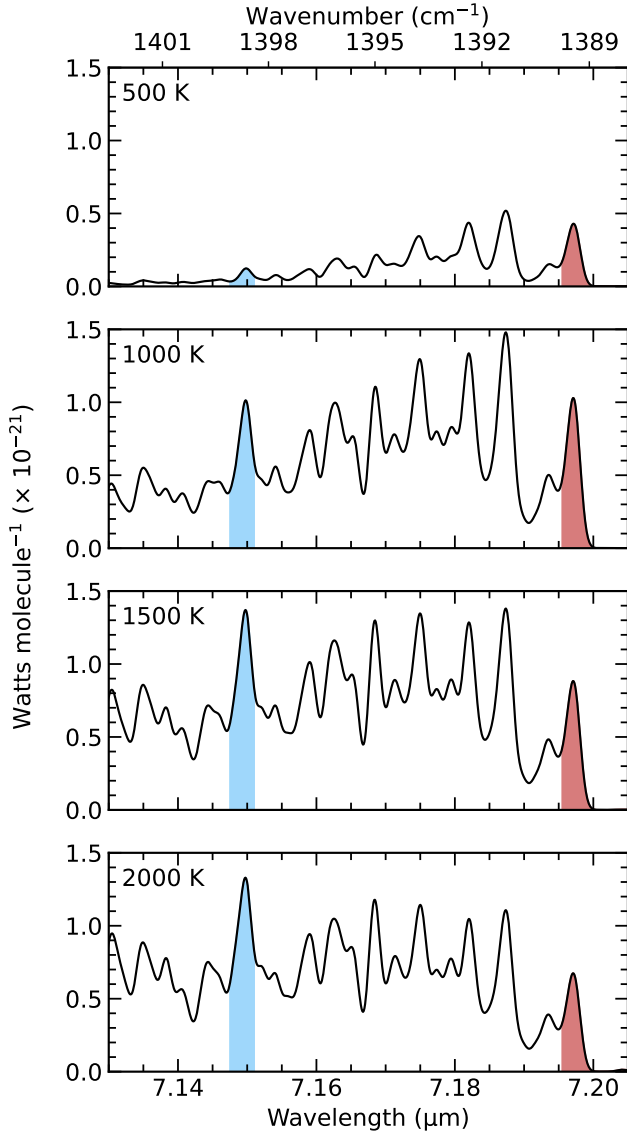


Figure 5. Model spectra of CH_3^+ at 500, 1000, 1500 and 2000 K. Bands at 7.15 and 7.2 μm are highlighted in blue and red, respectively.

4.2. Spectroscopic analysis of the CH_3^+ emission

CH_3^+ has a trigonal planar geometry (D_{3h} point group) and it lacks a permanent dipole moment. Because of that, it cannot be observed at radio wavelengths. However, it can be detected via ro-vibrational transitions at infrared wavelengths. CH_3^+ has symmetric (ν_1) and degenerate asymmetric (ν_3) stretching modes at around 3 μm ($\sim 3100 \text{ cm}^{-1}$), an out-of-plane umbrella bending mode (ν_2) and a degenerate in-plane (ν_4) bending mode at 7.2 μm ($\sim 1400 \text{ cm}^{-1}$; J. Meisner et al. 2019; P. B. Changala et al. 2023). Coriolis interactions between the ν_2 and ν_4 bands (also referred to as ν_2/ν_4 dyad) result in a characteristic pattern of ro-vibrational lines that occur in the 7.15–7.2 μm range.

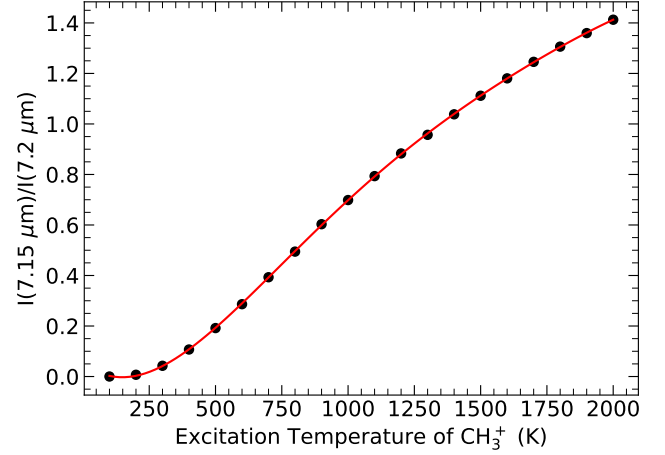


Figure 6. Intensity ratio $I(7.15\mu\text{m})/I(7.2\mu\text{m})$ as a function of excitation temperature of CH_3^+ . Black points represent the $I(7.15\mu\text{m})/I(7.2\mu\text{m})$ found in the model spectra for a given excitation temperature. A five-degree polynomial fit to those points is represented by the red curve. The polynomial is given by Eq. 1

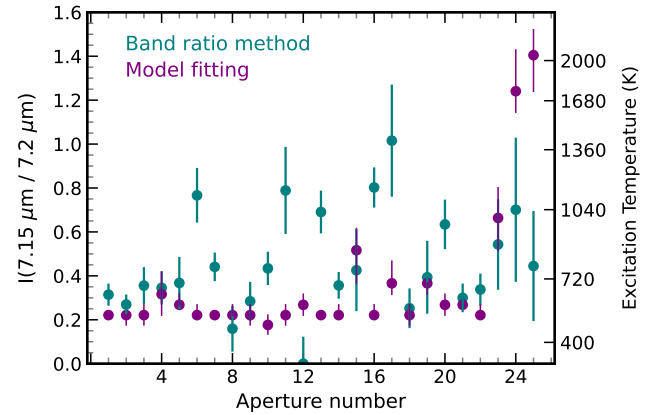


Figure 7. CH_3^+ excitation temperature derived for the 25 selected apertures. Teal points represent the $I(7.15\mu\text{m})/I(7.2\mu\text{m})$ band found in the observed extracted spectra. The corresponding temperatures (as described by Eq. 1) are shown on the y-axis on the right. Purple points represent the best-fit temperature derived from fitting models to the observations.

We compare the observations to the CH_3^+ model developed by P. B. Changala et al. (2023) using the latest and most detailed rovibrational assignments of the CH_3^+ feature based on theoretical and experimental studies. They performed detailed quantum mechanical calculations to derive spectroscopic constants (for the ν_2/ν_4 dyad) such as rotational constants, coriolis coupling constant, centrifugal distortion constants, etc., in the vibronic ground state of CH_3^+ . These constants were then validated by two different experimental methods. We use the PGOPHER (a general-purpose software for simulating electronic, vibrational and rotational spectra; C. M. Western 2017) file provided by (P. B. Changala et al.

2023) to compute the CH_3^+ model spectrum that can be compared to the observations at $7.2\mu\text{m}$ at JWST’s resolution. Note that the models assume a thermal distribution across rotational levels at $v=1$ level. We used PGOPHER to calculate CH_3^+ models spanning the temperature range from 100 K to 3000 K in steps of 50 K, where the spectrum is convolved with a Gaussian adopting a full width half maximum (FWHM) of 0.37 cm^{-1} , corresponding to the JWST resolving power of $R = 3600$ at $7.2\mu\text{m}$ (channel 1c; see I. Argyriou et al. 2023). We exported the models in the units of Watts per molecule.

Fig. 5 shows how the Q-branch of the CH_3^+ model spectrum changes as a function of excitation temperature. At JWST’s resolution, overlapping Q-branch lines cause a bump to appear. Note that the strength of different bands peaks at different temperatures based on the rotational levels of the transitions involved. For example, the strength of the $7.15\mu\text{m}$ band is lower than that of the $7.2\mu\text{m}$ band at lower temperatures. However, at temperatures above 1000K, the opposite is true. The intensity ratio of the $7.15\mu\text{m}$ band (shaded blue in Fig. 5) to the $7.20\mu\text{m}$ line (shaded red) is very small (<1) at low temperatures, ~ 1 at about 1300 K and much higher (>1) at high temperatures. This band ratio is thus a sensitive diagnostic of the CH_3^+ excitation temperature.

We will use this as a first estimate to determine the excitation temperature in the selected apertures, a diagnostic first used by M. Zannese et al. (2025) for the analysis of CH_3^+ emission in Orion Bar PDR. We fit a gaussian with a linear baseline to the $7.15\mu\text{m}$ ($7.147 - 7.151\mu\text{m}$) band and $7.20\mu\text{m}$ ($7.195 - 7.199\mu\text{m}$) band, for both models and observations. The ratio of $I(7.15\mu\text{m})/I(7.2\mu\text{m})$ derived from the model spectra versus excitation temperature is shown in Fig. 6, which exhibits the behaviour described above. This ratio should change gradually as a function of temperature, and to obtain a continuous form of this function, we fitted a polynomial to the data points. We found that a fifth-order polynomial fits the individual points well (see Fig. 6) using the following coefficients:

$$T(x) = 1883x^5 - 7260x^4 + 10640x^3 - 6956x^2 + 2941x + 159.6 \quad (1)$$

where x is the band ratio. We can then use this polynomial to determine the excitation temperature using measured band ratios in the observational spectra. For each of the 25 apertures, we applied the same fitting procedure to the $7.15\mu\text{m}$ and $7.20\mu\text{m}$ bands as done for the models. These fits are shown in Appendix C, and the intensity and uncertainties on the bands and the band ratios are listed in Table 1. The uncertainties in the observations determine uncertainties in the intensity. The teal points in Fig. 7 show the band ratio and temperatures derived for each aperture based on Eq. 1. All apertures exhibit excitation temperatures of 500–1000 K (band

ratios of 0.15–0.80). While the reported uncertainties primarily reflect random errors, systematic effects—especially those introduced by the linear baseline fit—are harder to quantify and could artificially increase or decrease the band ratios in certain apertures.

We can also determine the excitation temperatures by directly fitting model spectra to the continuum-subtracted spectra, with the temperature and column density of CH_3^+ as free parameters. We used the continuum-subtracted flux (in $\text{W m}^{-2}\mu\text{m}^{-1}$) in the range of $7.13\text{--}7.199\mu\text{m}$ to fit models to observations. We used the `lmfit` (M. Newville et al. 2014) python package, adopting a Levenberg-Marquardt algorithm to determine temperatures and column densities; the excitation temperatures, however, were constrained on a grid with values between 100 K and 3000 K in steps of 50 K. We also allowed for small wavelength shifts ($\sim \pm 0.0015\mu\text{m}$) during fitting to account for radial velocity corrections. Appendix D shows the resulting best fits for all 25 apertures, and the resulting best fit parameters are listed in Table 1. Parameter uncertainties presented in Table 1 denote 1σ confidence intervals, obtained from the projections of the $\Delta\chi^2 = 1$ contour on the χ^2 surface. The models generally reproduce the observations extremely well, leaving only very small residuals in this wavelength range. We find that the excitation temperature is in the range of 500–800K for apertures 1-22 and temperature 1000-2000K for apertures 23-25; these are represented by purple points in Fig 7.

Although no clear patterns exist in excitation temperature variations across the inner bubble and torus, the outer bubble exhibits significantly higher temperatures (1000–2000 K) compared to 500–800 K elsewhere. The column densities (Table 1) are typically a few times 10^{12} cm^{-2} , ranging about two orders of magnitude from the lowest (aperture 23) to highest (aperture 1) values. The highest column density is seen where the inner bubble runs into the rear side of the torus. These column densities are however, not an accurate representation of the total CH_3^+ column density: as was pointed out by O. Berné et al. (2023) and M. Zannese et al. (2025) since the CH_3^+ emission originates from chemical pumping and consequently, the excitation of the vibrational levels is far from LTE.

5. DISCUSSION

5.1. Formation of CH_3^+ in NGC 6302

The detection of CH_3^+ in NGC 6302 was surprising at first because this is an O-rich PN where most carbon is locked up in CO, making carbon chemistry unlikely. However, intense UV radiation from the central star fundamentally changes this scenario by providing two key ingredients required for the formation of CH_3^+ : C^+ ions and vibrationally excited H_2 . C^+ reacts with the vibrationally excited H_2 to form CH^+ , followed by two successive exothermic hydrogen abstraction

Table 1. The excitation temperature, column density and wavelength shift for each of the 25 apertures from Fig. 2.

Aperture	$I_{7.15} \pm \sigma$ [10^{-1}]	$I_{7.20} \pm \sigma$ [10^{-1}]	Ratio $\pm \sigma$ [10^{-1}]	T^a [K]	T^b [K]	$N \pm \sigma_N$ [10^{12} cm^{-2}]	wavelength shift [$10^{-3} \mu\text{m}$]
1	8.3 ± 1.3	26.4 ± 1.3	3.1 ± 0.5	660^{+40}_{-50}	550^{+0}_{-0}	13.3 ± 0.2	-1.02 ± 0.03
2	7.2 ± 1.2	26.6 ± 1.2	2.7 ± 0.5	620^{+40}_{-50}	550^{+0}_{-50}	11.4 ± 0.2	-0.97 ± 0.04
3	4.0 ± 0.9	11.2 ± 0.9	3.6 ± 0.8	700^{+70}_{-80}	550^{+50}_{-50}	3.8 ± 0.2	-0.84 ± 0.09
4	5.5 ± 1.1	16.0 ± 1.1	3.5 ± 0.8	690^{+70}_{-70}	650^{+100}_{-100}	2.0 ± 0.1	-1.17 ± 0.1
5	4.3 ± 1.3	11.8 ± 1.3	3.7 ± 1.2	710^{+100}_{-110}	600^{+50}_{-0}	5.4 ± 0.1	-1.18 ± 0.05
6	8.9 ± 1.2	11.7 ± 1.2	7.7 ± 1.2	1110^{+150}_{-140}	550^{+50}_{-0}	5.8 ± 0.2	-1.19 ± 0.06
7	6.7 ± 0.9	15.1 ± 0.9	4.4 ± 0.7	770^{+60}_{-60}	550^{+0}_{-0}	6.9 ± 0.1	-1.16 ± 0.04
8	1.4 ± 0.9	8.6 ± 0.9	1.6 ± 1.0	490^{+120}_{-190}	550^{+50}_{-0}	4.5 ± 0.1	-1.19 ± 0.06
9	3.5 ± 1.0	12.3 ± 1.0	2.8 ± 0.9	630^{+80}_{-100}	550^{+50}_{-50}	4.1 ± 0.2	-1.12 ± 0.1
10	7.7 ± 1.2	17.7 ± 1.2	4.3 ± 0.8	770^{+70}_{-70}	500^{+50}_{-50}	5.8 ± 0.2	-1.12 ± 0.06
11	4.8 ± 1.0	6.1 ± 1.0	7.9 ± 2.0	1140^{+250}_{-220}	550^{+50}_{-50}	3.3 ± 0.2	-0.9 ± 0.1
12	0.0 ± 1.7	13.7 ± 1.7	0.0 ± 1.2	160^{+270}_{-490}	600^{+50}_{-50}	5.7 ± 0.1	-1.01 ± 0.05
13	8.0 ± 0.9	11.6 ± 0.9	6.9 ± 1.0	1020^{+110}_{-110}	550^{+0}_{-0}	5.4 ± 0.1	-0.95 ± 0.05
14	3.7 ± 0.6	10.4 ± 0.6	3.6 ± 0.6	700^{+50}_{-50}	550^{+50}_{-0}	4.5 ± 0.1	-1.1 ± 0.07
15	2.3 ± 0.9	5.4 ± 0.9	4.3 ± 1.9	760^{+180}_{-170}	850^{+100}_{-150}	1.0 ± 0.1	-0.74 ± 0.13
16	10.1 ± 0.9	12.5 ± 0.9	8.0 ± 0.9	1150^{+120}_{-110}	550^{+50}_{-0}	4.6 ± 0.2	-1.14 ± 0.07
17	4.2 ± 0.7	4.1 ± 0.7	10.2 ± 2.6	1430^{+380}_{-320}	700^{+100}_{-50}	1.4 ± 0.1	-1.13 ± 0.12
18	3.4 ± 1.2	13.6 ± 1.2	2.5 ± 0.9	600^{+90}_{-110}	550^{+50}_{-50}	4.7 ± 0.2	-1.25 ± 0.07
19	1.7 ± 0.7	4.2 ± 0.7	3.9 ± 1.7	730^{+150}_{-160}	700^{+0}_{-50}	1.8 ± 0.1	-1.34 ± 0.09
20	5.4 ± 0.8	8.6 ± 0.8	6.3 ± 1.1	960^{+130}_{-110}	600^{+50}_{-50}	2.2 ± 0.1	-1.16 ± 0.08
21	3.2 ± 0.7	10.8 ± 0.7	3.0 ± 0.7	650^{+60}_{-70}	600^{+0}_{-0}	4.8 ± 0.1	-0.81 ± 0.05
22	4.6 ± 0.9	13.5 ± 0.9	3.4 ± 0.7	680^{+60}_{-70}	550^{+50}_{-0}	4.8 ± 0.1	-0.85 ± 0.05
23	0.7 ± 0.2	1.4 ± 0.2	5.4 ± 2.1	870^{+220}_{-180}	1000^{+150}_{-150}	0.2 ± 0.0	-0.96 ± 0.11
24	1.0 ± 0.4	1.4 ± 0.4	7.0 ± 3.3	1030^{+410}_{-320}	1750^{+350}_{-150}	0.4 ± 0.0	-1.05 ± 0.1
25	0.9 ± 0.5	2.0 ± 0.5	4.4 ± 2.5	780^{+250}_{-240}	2050^{+500}_{-300}	0.4 ± 0.0	-1.05 ± 0.11

^aBand ratio method

^bModel fitting

reactions to produce CH_3^+ (O. Berné et al. 2023; M. Zan-
nese et al. 2025). The UV radiation plays a crucial role be-
cause it dissociates and photoionizes CO, thereby providing
 C^+ . Moreover, in such environments, the energy barrier of
the CH^+ step ($\approx 0.40 \text{ eV}$ or $\sim 4640 \text{ K}$) is overcome by FUV-
pumped H_2 , which leaves H_2 in a vibrationally excited state
(M. Agúndez et al. 2010; Z. Nagy et al. 2013; B. Godard &
J. Cernicharo 2013). CH^+ and CH_3^+ form in a vibrationally
excited state and then relax radiatively. This excitation mech-
anism is called chemical (formation) pumping, and the ob-
served emission thus traces the newly formed CH^+ and CH_3^+ .

Our observations confirm this formation pathway in
NGC 6302 through spatial comparison of ^{12}CO , H_2 and CH_3^+ .
Fig. 8 shows that ^{12}CO and H_2 are indeed present where CH_3^+
emission is detected i.e. where CH_3^+ is formed. CH_3^+ and

^{12}CO are also present in the outer bubble, but as mentioned
above, their column densities are too low to be visible in the
integrated surface brightness map. Since the vibrationally ex-
cited H_2 lines detected with JWST/MIRI data have very low
S/N, their spatial distribution has large uncertainties for com-
parison with CH_3^+ distribution. We therefore use the H_2 0-0
S(5) line (at $6.908 \mu\text{m}$) as a tracer for the H_2 distribution.

Fig. 9 shows the normalized surface brightness profile of
these species along the cut indicated in Fig. 2, zoomed in to
show the east and west edges of the bubble. On the east edge,
 CH_3^+ , ^{12}CO and H_2 show peaks in the surface brightness pro-
files at the same distance from the star, with CH_3^+ and H_2
having very similar profiles. The west edge shows a very dif-
ferent distribution, likely because of projection effects. No-
tably, CH_3^+ exhibits the sharpest rise and fall at the inner bub-
ble edge compared to other species, including H_2 . This could

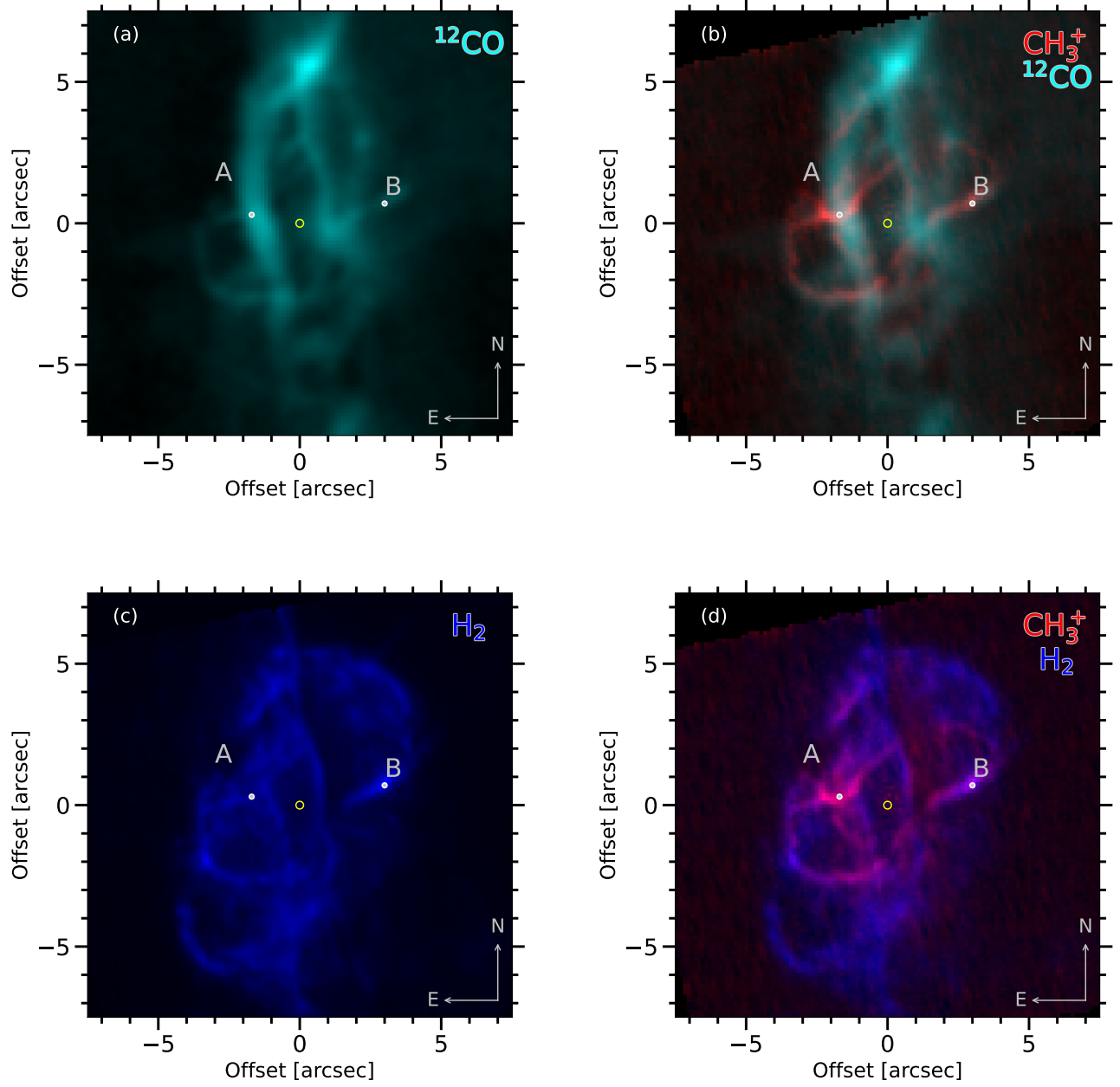


Figure 8. Comparison of the integrated surface brightness maps of CH_3^+ , H_2 and ^{12}CO . (a) ^{12}CO (2–1, 230GHz, by Moraga Baez et al., 2025, under review). (b) Two colour composite image: CH_3^+ (red) and ^{12}CO (cyan). (c) H_2 0-0 S5 at $6.908\ \mu\text{m}$ (d) Two colour composite image: CH_3^+ (red) and ^{12}CO (cyan). Locations marked by A and B represent the two bright spots on CH_3^+ map. The maps are centred at the central star (R.A.= $17^{\text{h}}:13^{\text{m}};44.488 \pm 0.004^{\text{s}}$, Dec. = $-37^{\circ}:06':11.76 \pm 0.03'$) whose position is indicated by the yellow circle at the center.

indicate where FUV-pumped H_2 ($v>0$) is available; however, NIRSPEC observations of vibrationally excited H_2 lines are needed for confirmation. Alternatively, this may reflect the influence of fast winds or energetic shocks that formed the inner bubble (Matsuura et al., 2025, MNRAS, in press), confining CH_3^+ formation to a narrow region along the edge.

5.2. Spatial variations in CH_3^+ emission

CH_3^+ emission shows significant spatial variations in column density across NGC 6302. The most prominent features are two bright spots, Points A and B, which exhibit similar temperatures to other regions but significantly higher column densities (See Fig. 2). This enhancement occurs where the inner bubble runs into the torus, giving the inner bubble

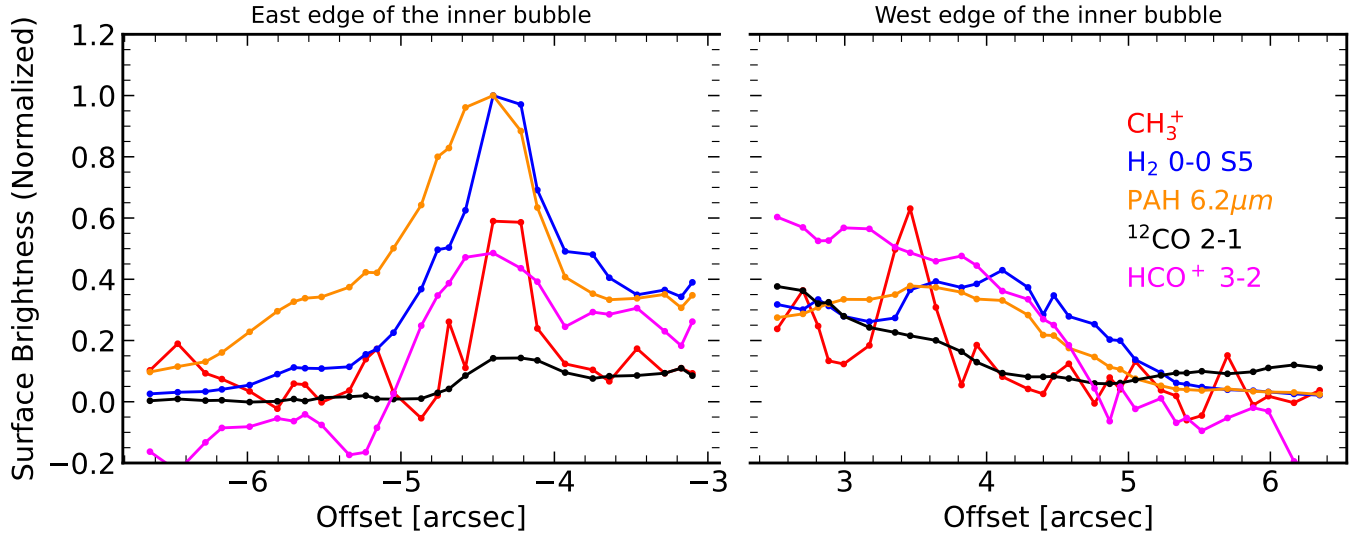


Figure 9. Normalized surface brightness profiles of CH_3^+ , H_2 , PAH $6.2\ \mu\text{m}$ (by Clark et al., in prep), CO and HCO^+ (by Moraga Baez et al., 2025, under review) along the cut. Offset is measured in arcsec from the position of the central star (R.A. = $17^{\text{h}}:13^{\text{m}};44.488 \pm 0.004^{\text{s}}$, Dec. = $-37^{\circ}:06':11.76 \pm 0.03'$).

its characteristic peanut shape and creating regions of higher gas density. Point A (region around apertures 1,2, and 12), exhibiting the brightest CH_3^+ emission, is located at the intersection of the inner bubble with the rear side of the torus. A corresponding bright feature is not observed where the inner bubble intersects the front side because of the extinction by the dust in the torus. However, enhanced emission is detected in both the east (aperture 7) and west (apertures 5 and 6) of this intersection, potentially tracing redirected material emerging from the intersection. This spot on the west is designated point B. Apart from higher gas density, secondary factors may also contribute to enhanced formation rates. For example, at point A, the H I emission is several times brighter than the surrounding regions, suggesting higher UV flux, and point B coincides with a brighter H_2 emission that could boost CH_3^+ production.

In contrast, the outer bubble shows column densities an order of magnitude lower than typical values in the inner bubble and the torus and two orders of magnitude lower than point A. This is likely due to the lower column density of CO in the outer bubble. Interestingly, the CH_3^+ excitation temperature is significantly higher in the outer bubble (1000-2000K) than what is found in the inner bubble and torus (500-800K). The chemical pumping model by M. Zannese et al. (2025) shows that higher excitation temperature of CH^+ results from enhanced H_2 level population densities, which occur when H_2 is more efficiently FUV-pumped to higher vibrational states. Similarly, hotter CH_3^+ can form if the H_2 involved in hydrogen abstraction reactions is vibrationally excited to higher levels. This indicates that the outer bubble is less shielded than the inner structures.

5.3. Chemistry initiated by CH_3^+

The main takeaway from the CH_3^+ detection in NGC 6302 is recognizing that hydrocarbon radical chemistry, previously overlooked in oxygen-rich environments, should be incorporated into planetary nebula chemical models. CH_3^+ drives gas-phase carbon chemistry through exothermic ion-neutral reactions, with models showing it can facilitate the formation of numerous species, including CN, HCN, CH_2CO^+ , CH_3OH , H_2CO , $\text{C}_2\text{H}_5\text{OH}$, CH_3OCH_3 , H_2CN^+ , etc. (see Fig. 2 by D. Smith (1992) for the summary of the chemical network initiated by CH_3^+).

Spatial coincidences in NGC 6302 suggest potential connections between CH_3^+ and other observed species that warrant consideration in chemical models. CH_3^+ is spatially coincident with HCO^+ along the inner and outer bubble edges (see Fig. 10), consistent with the exothermic reaction $\text{CH}_3^+ + \text{O} \rightarrow \text{HCO}^+ + \text{H}_2$ (G. B. I. Scott et al. 2000). The co-location of CH_3^+ with PAHs (see Fig. 11) raises the question whether CH^+ or CH_3^+ might initiate chemistry relevant to PAH formation. M. Agúndez et al. (2008) reported that strong FUV/X-ray radiation can efficiently drive carbon chemistry in dense O-rich gas where CH^+ and CH_3^+ can further react with e^- , C, and H_2 to form C_2H_2 , which can then initiate the chemistry leading to aromatic species like PAHs if activation energies are overcome. Since PAH emission is widespread in NGC 6302's central region—including the torus, inner and outer bubbles where CH_3^+ is detected—the role of CH_3^+ in the bottom-up PAH formation sequence should be investigated.

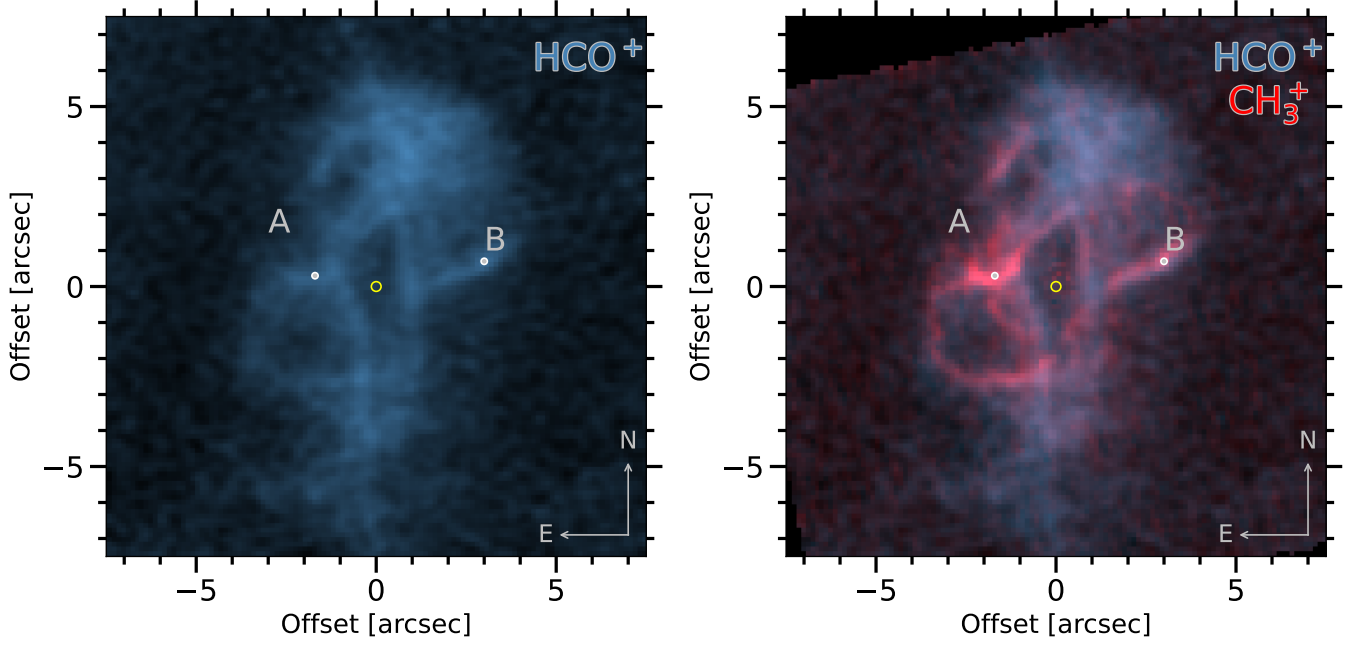


Figure 10. HCO^+ map (by Moraga Baez et al., 2025, under review) is shown in blue on the left and an overlay of CH_3^+ map is shown in red on the right. Locations marked by A and B represent the two bright spots on CH_3^+ map. The maps are centred at the central star (RA=17:13:44.488 $\pm$ 0.004, dec=-37:06:11.76 $\pm$ 0.03) whose position is indicated by the white circle at the center.

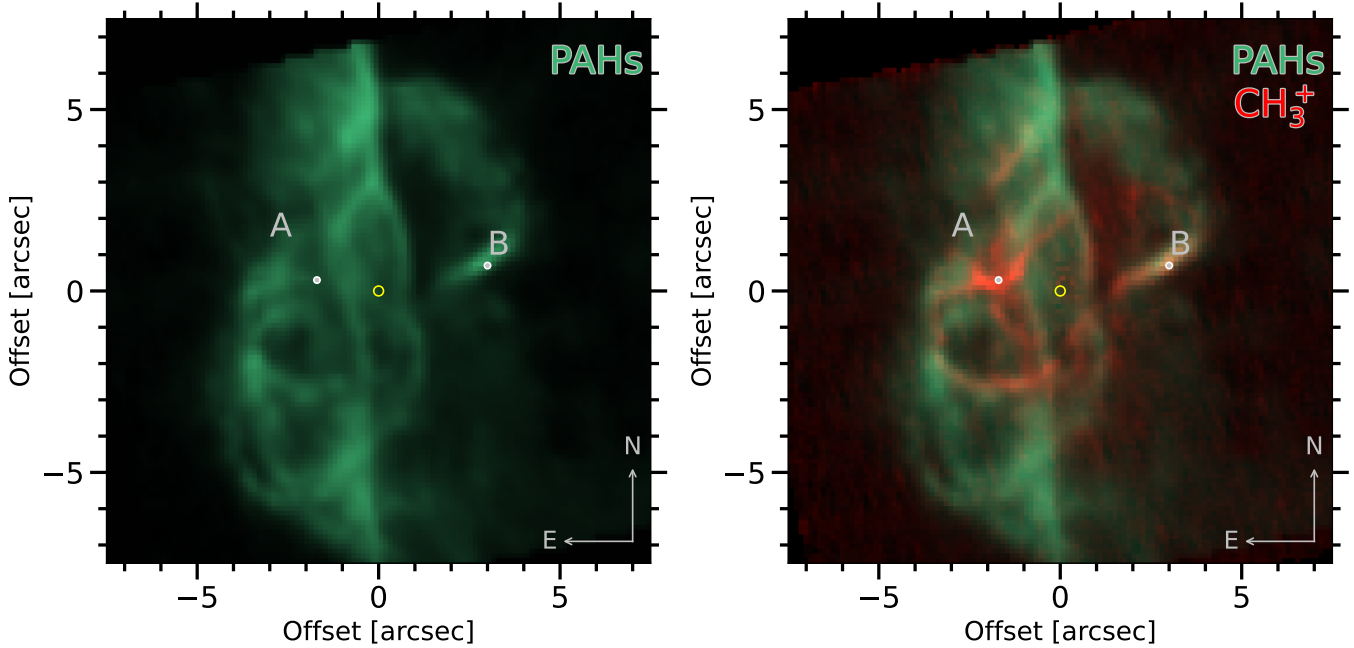


Figure 11. The PAH map (by Clark et al., in prep) here on the left shows the integrated intensity of the 6.2 μm feature (integrated over the 6.12–6.5 μm range after continuum subtraction). A comparison with CH_3^+ map is shown on the right. Locations marked by A and B represent the two bright spots on CH_3^+ map. The maps are centred at the central star (R.A.= 17^h:13^m:44.488 $\pm$ 0.004^s, Dec. = -37^o:06':11.76 $\pm$ 0.03[']) whose position is indicated by the white circle at the center.

6. SUMMARY & CONCLUSIONS

We report the detection of methyl cation (CH_3^+) in JWST-MIRI observations of the O-rich planetary nebula NGC 6302,

marking the first identification of this key organic precursor in such an environment. We investigated the nature of the CH_3^+ emission and its relation to other species such as ^{12}CO ,

H_2 , PAHs and HCO^+ . We fit the observed CH_3^+ emission with LTE models to derive excitation temperature and column density. Model fits reproduce the CH_3^+ emission very well, yielding excitation temperatures of 500-800K in the inner bubble and the torus, **significantly higher excitation temperature of 1000-2000K in the outer bubble**. Column densities of CH_3^+ in the excited state are typically $\sim 10^{12} \text{ cm}^{-2}$ ranging from $\sim 10^{11} \text{ cm}^{-2}$ **in the outer bubble** to $\sim 10^{13} \text{ cm}^{-2}$ where the inner bubble runs into the torus. **These measurements represent only a fraction of the total population due to non-LTE chemical pumping. The spatial coincidence of CH_3^+ with ^{12}CO and H_2 supports its formation via chemical pumping in the UV-irradiated O-rich gas.**

The presence of CH_3^+ in this O-rich PN highlights the carbon chemistry at play in O-rich environments under strong UV irradiation. These results indicate that hydrocarbon radical chemistry must be incorporated into chemical models of PNe. Further observations at comparable spatial resolution, particularly in the near-IR (which is rich in molecular features), are essential to unravel the different chemical networks involved in this chemistry.

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APPENDIX

A. CONTINUUM

An example of the continuum adopted for fitting the CH_3^+ spectral features is shown in Fig. 12.

B. H_2 MAP WITH APERTURES

The 25 apertures used to sample the CH_3^+ emission across distinct regions are plotted on the H_2 map in Fig. 13.

C. BAND RATIO METHOD

The best-fit Gaussians to the $7.15\mu\text{m}$ and $7.2\mu\text{m}$ bands for all 25 apertures are shown in Figs. 14 and 15, respectively. These are used to determine the excitation temperature of CH_3^+ using the band ratio method as described in Sec. 4.

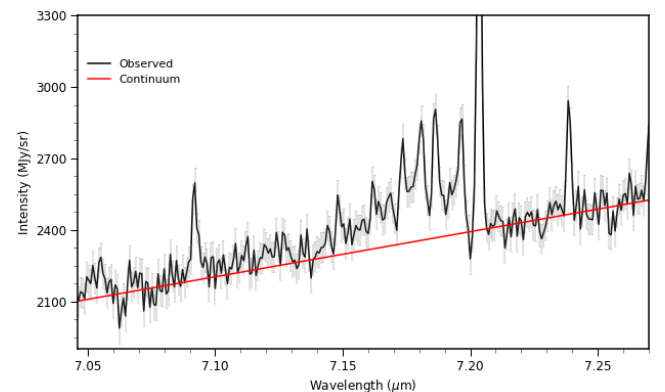


Figure 12. JWST MIRI spectrum at aperture 1 (in black). The red line provides an example of the continuum adopted to fit the models to the observations.

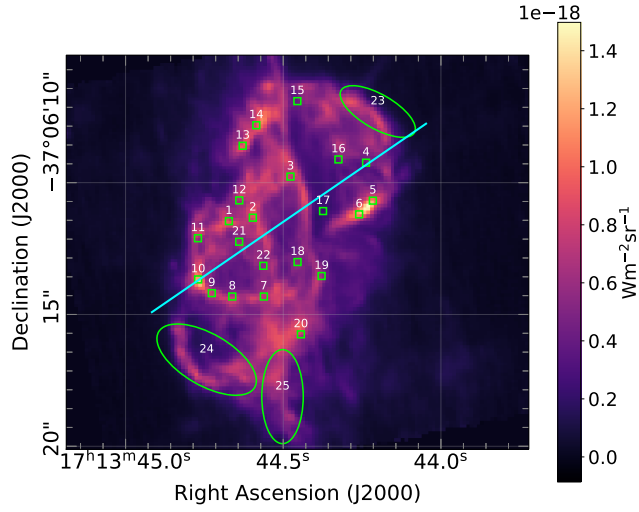


Figure 13. Integrated surface brightness map of H₂ 0-0 S(5) line at 6.908 μ m. The green boxes and ellipses represent the 25 apertures used to study CH₃⁺ emission in detail.

D. FIT RESULTS

The best-fit models for determining CH₃⁺ temperature and column density in all 25 apertures are shown in Fig. 16. See Sec. 4 for details on the fitting procedure.

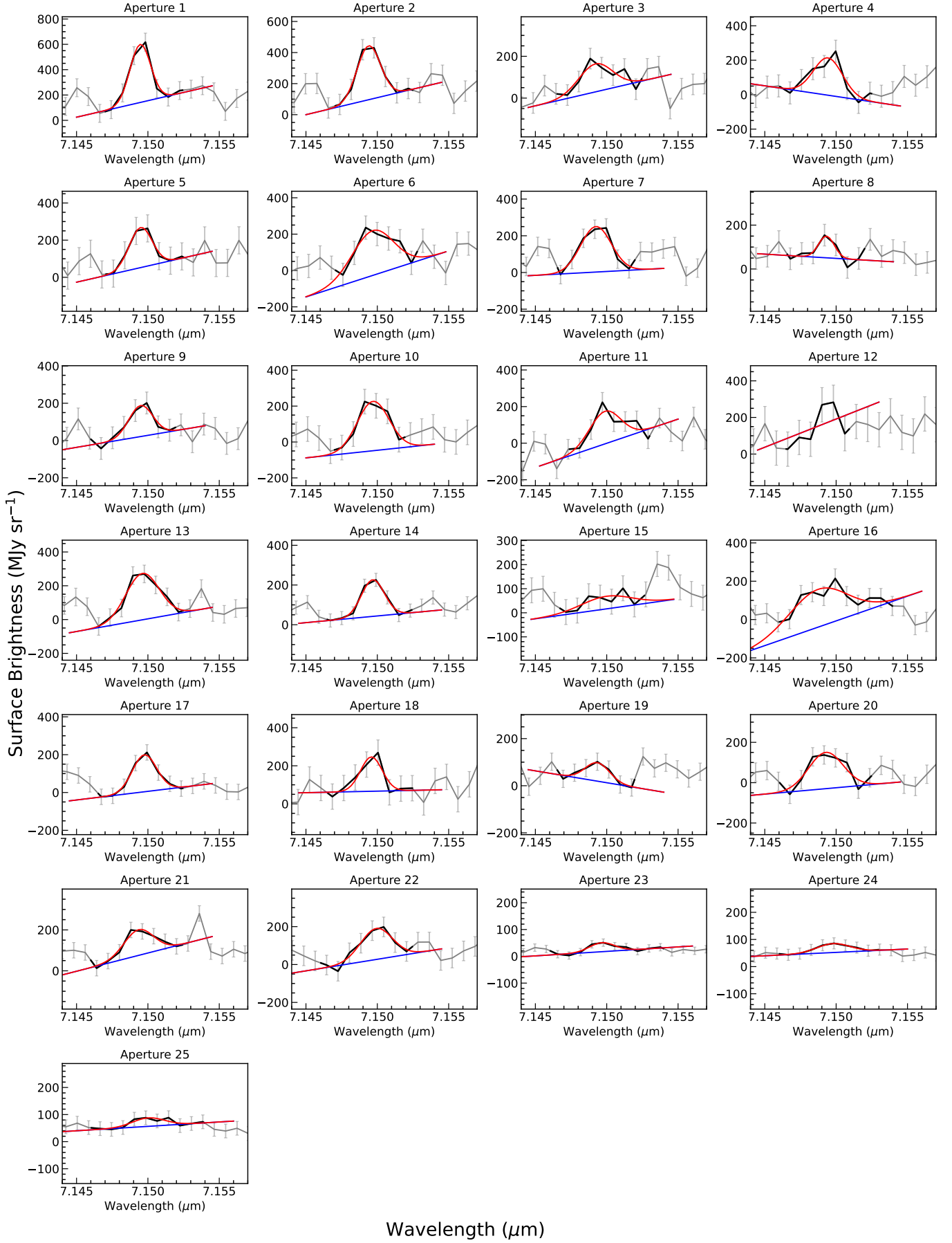


Figure 14. Gaussian fits to the 7.15 band used for the band ratio method to determine the temperature. The intensity of the band at each aperture is listed in Table 1.

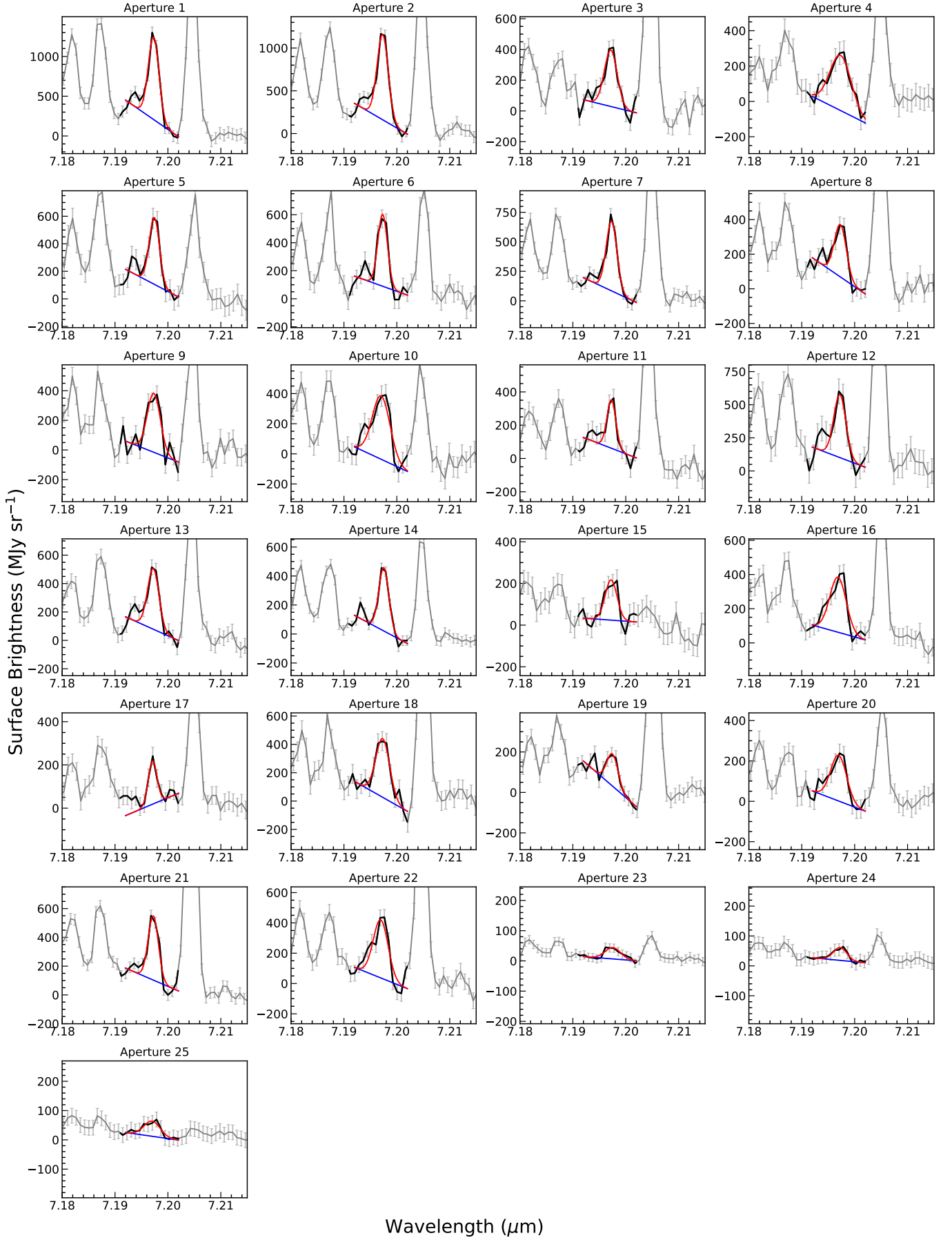
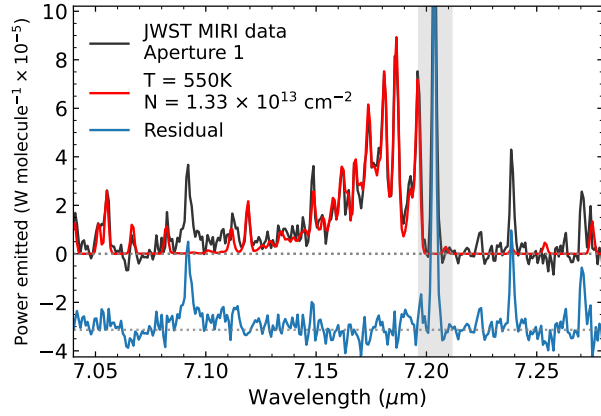
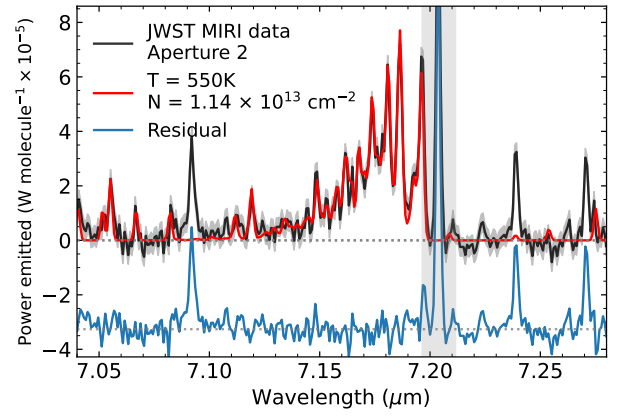


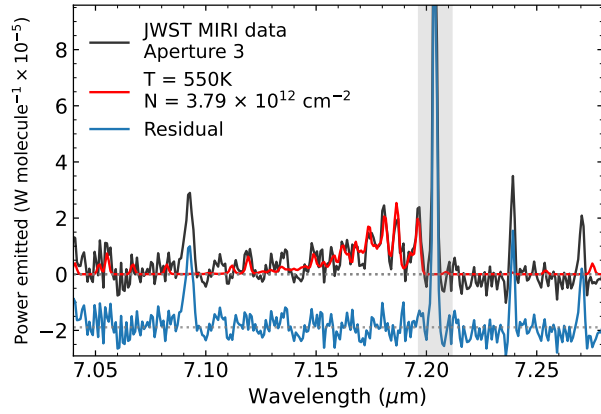
Figure 15. Gaussian fits to the 7.2 band used for the band ratio method to determine the temperature. The intensity of the band at each aperture is listed in Table 1.



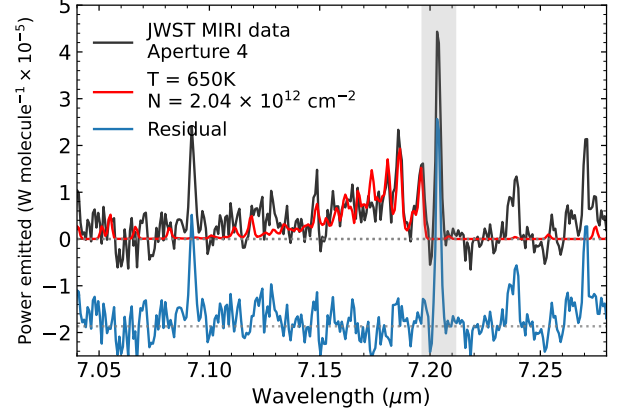
(a) Aperture 1.



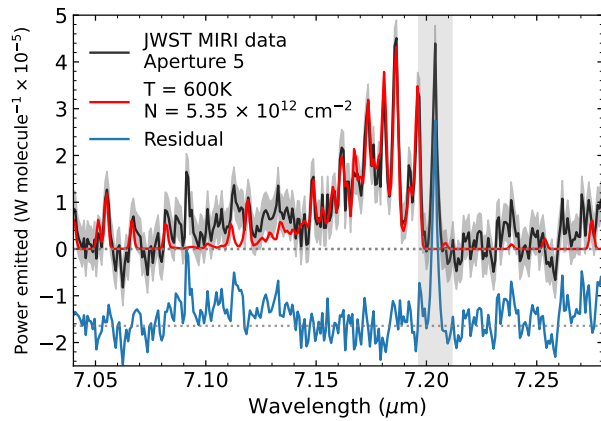
(b) Aperture 2.



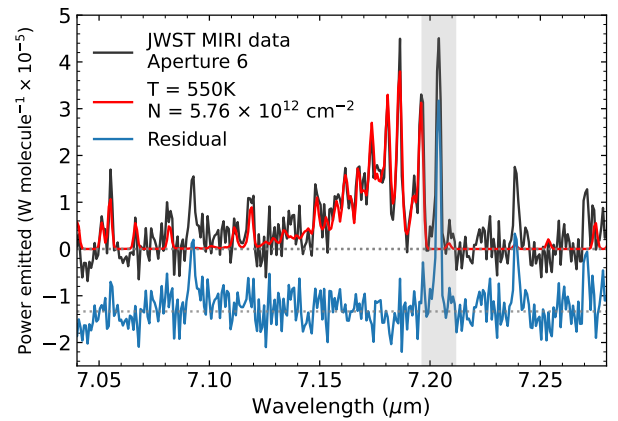
(c) Aperture 3.



(d) Aperture 4.

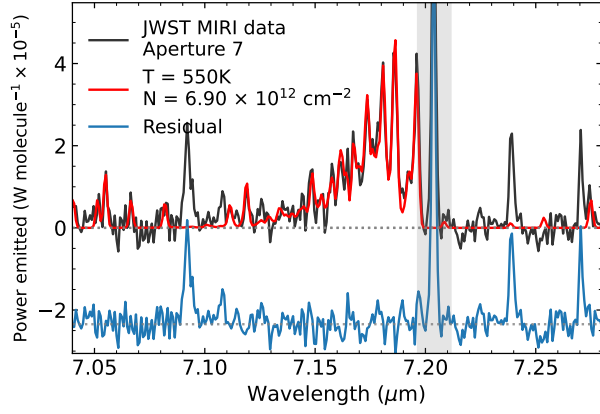


(e) Aperture 5.

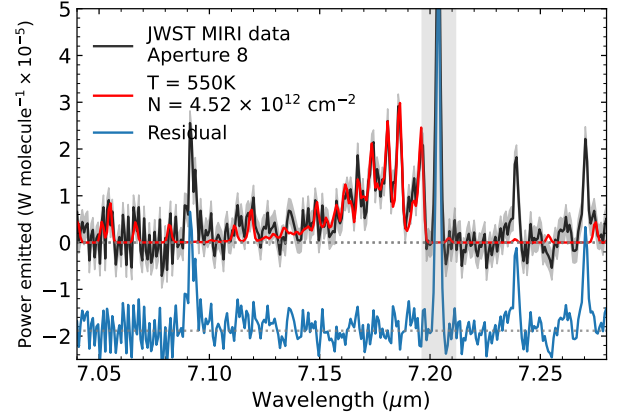


(f) Aperture 6.

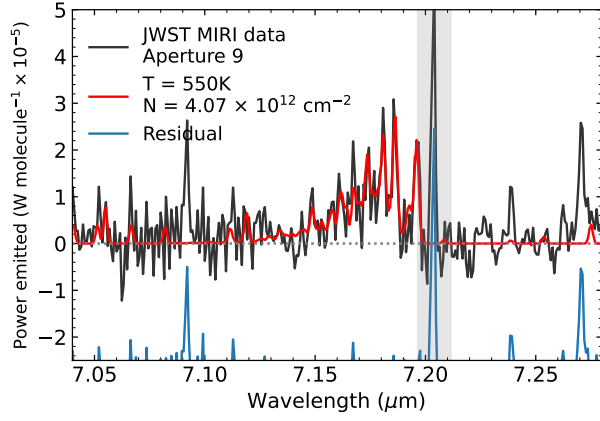
Figure 16. Observations (black), best-fit model (red), and residuals (blue) for each aperture.



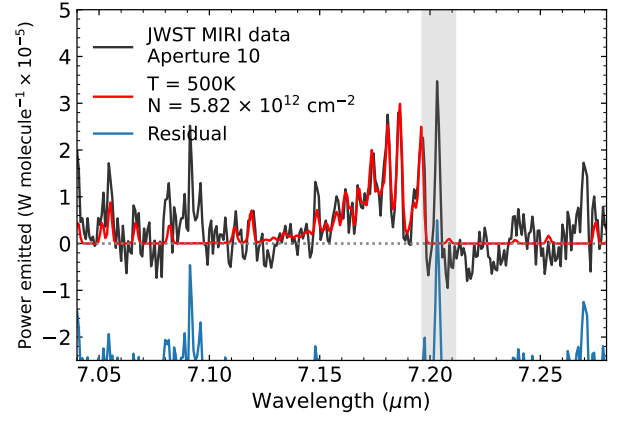
(g) Aperture 7.



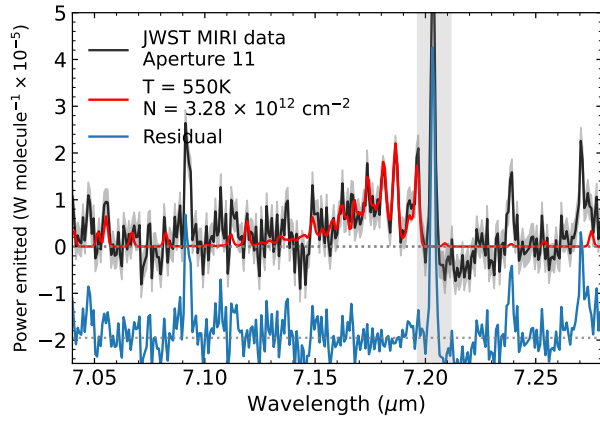
(h) Aperture 8.



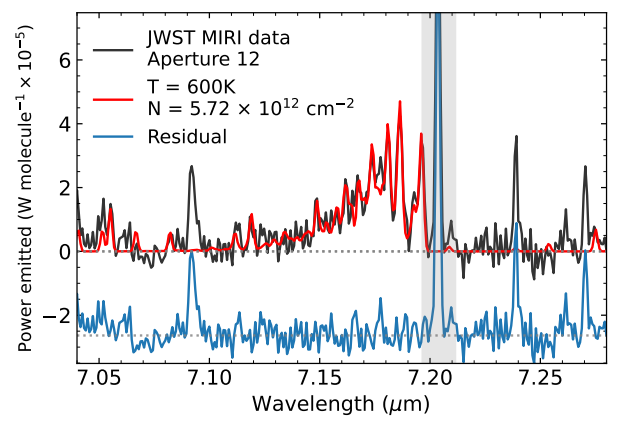
(i) Aperture 9.



(j) Aperture 10.

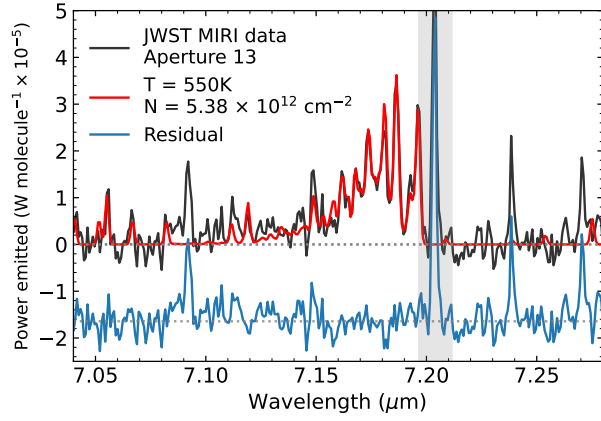


(k) Aperture 11.

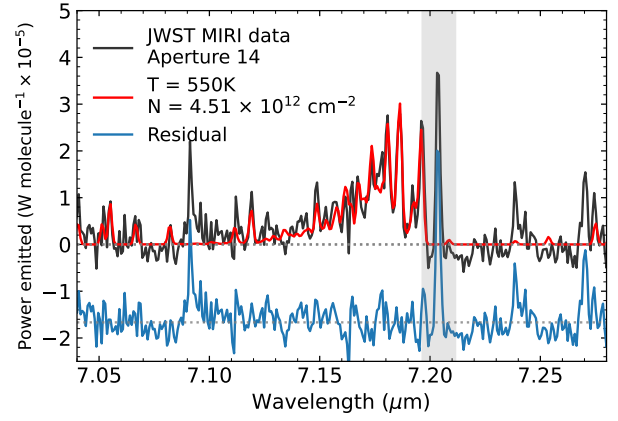


(l) Aperture 12.

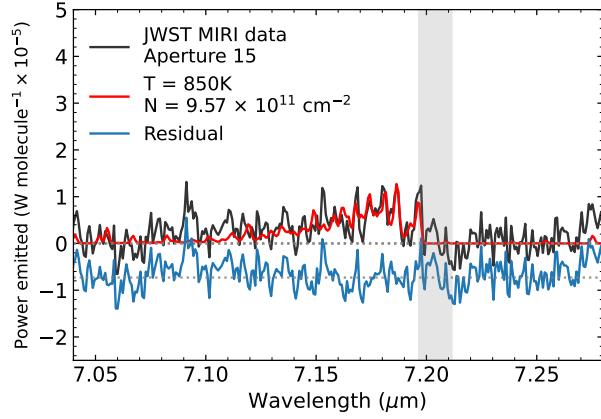
Figure 16. Figure 16 (continued).



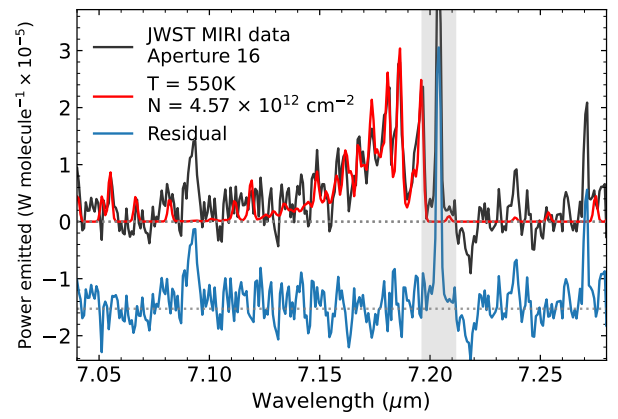
(m) Aperture 13.



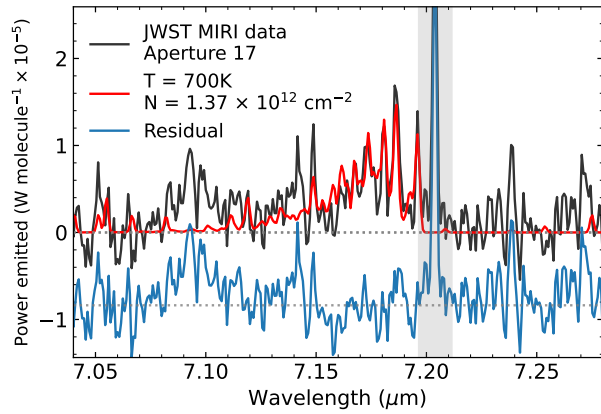
(n) Aperture 14.



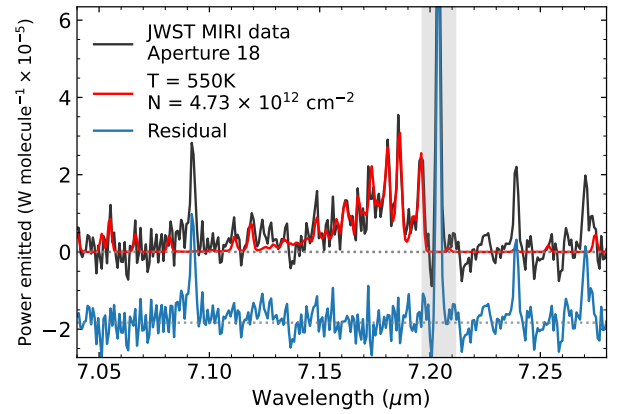
(o) Aperture 15.



(p) Aperture 16.

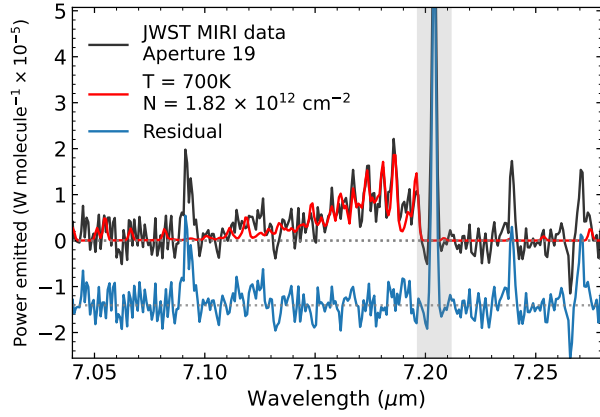


(q) Aperture 17.

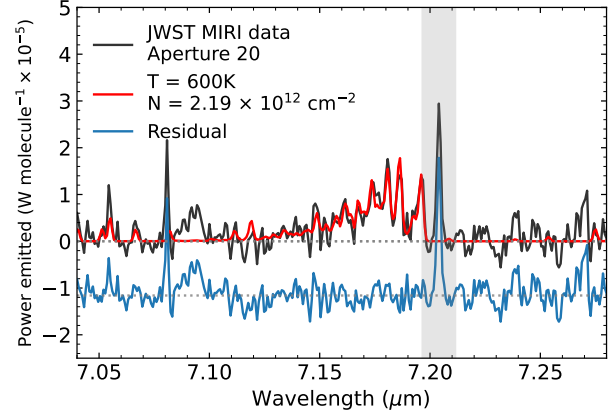


(r) Aperture 18.

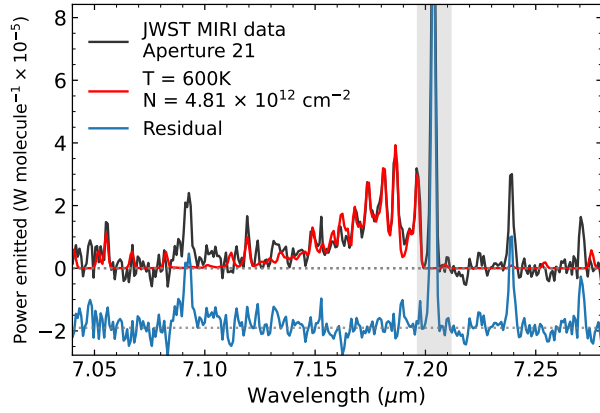
Figure 16. Figure 16 (continued).



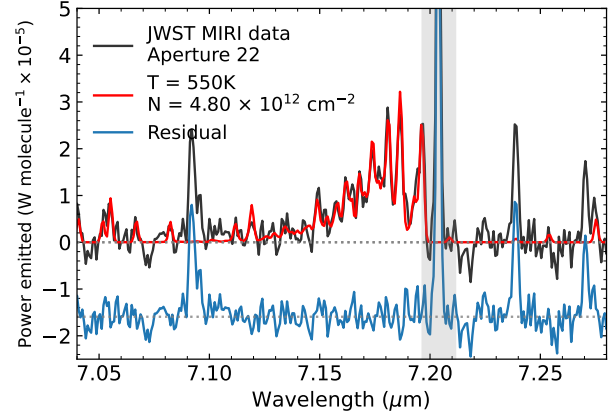
(s) Aperture 19.



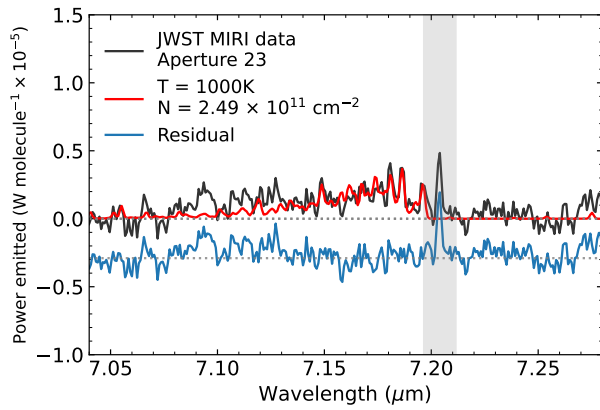
(t) Aperture 20.



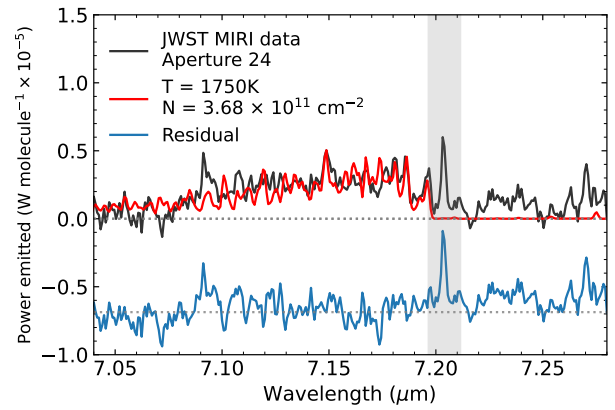
(u) Aperture 21.



(v) Aperture 22.

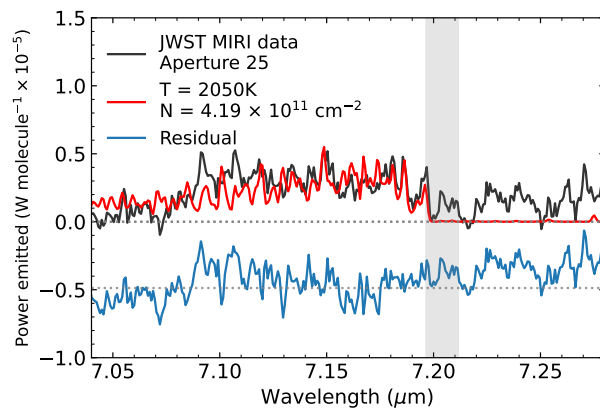


(u) Aperture 23.



(v) Aperture 24.

Figure 16. Figure 16 (continued).



(s) Aperture 25.

Figure 16. Figure 16 (continued).

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