Unveiling the origin and excitation mechanisms of the warm CO, OH and CH⁺



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Observations

- Fully sampled observations with PACS: 4×4 rastermap (OH 85 μm: 5x5) covering an area of ~100" x 100". Each position is composed of 5 × 5 spatial pixels (spaxel). For the spaxels, the line is observed in 25 positions

- Spatial resolution between 6" to 10" (0.01-0.02 pc)

- Targeted lines: CO |=19-18 (137μm), CH⁺ |=3-2 (cube includes: CO |=22-21, OH 119μm), OH 84µm, [OI] 63µm, [OI] 145µm (cube includes: CO J=18-17), and [CII] 158 µm (Bernard-Salas et al. 2012)

- PACS data was processed with HIPE 10 standard pipeline, and further with PACSman (Lebouteiller et al. 2012)

- SPIRE observations on ¹²CO (transitions: J=5-4 to J=13-12) and ¹³CO (transitions: J=5-4 to J=11-10). The data was processed with HIPE 11 standard pipeline. Supreme method (see poster by H.Ayasso, Köhler et al. (submitted)) also used with HIPE 8 as HIPE 11 version is still under development. This method resolves the structures in the bar with more detail than the standard pipeline (see next section).

Spatial distribution and excitation of the warm CO

The morphologies of CO J=19-18 and the ¹³CO J=11-10 are very similar, only shifted as the ¹³CO is a lower transition

The edge of CO J=19-18 agrees well with rovibrationally excited H₂ (H₂^{*}) which is very sensitive to UV field, we thus conclude that high-J CO directly traces the irradiated dense structures.

The excitation temperature is high where the PAH (poly-aromatic hydrocarbons) emission peaks due to UV heating. This puts strong constraints on the origin of the CO excitation: UV heating dominates

CH⁺ and OH comparison with H₂

CH⁺ J= 3-2 and OH 85 μ m have similar critical densities (~10¹⁰ cm⁻³) and upper energy levels (~250 K) which make them ideal for comparison



Contours of ¹³CO J=11-10 (Supreme method) and H₂ 1-0 S(1) line (Walmsley et al. 2000). The clumps are seen in the high-J ¹²CO and intermediate-J ¹³CO (see the arrows).



Contours of CO |=19-18.

The ratio peaks in front of the bar.



Contours of ratio: CO J=19-18 / CO J=11-10

CH⁺ formation: CH⁺ destruction:

Map: H2 I-0 S(I) CH⁺ contour levels at 0.3, 0.5, 0.7, and 0.9 of the emission maximum

CH⁺ comparison with H₂:

 $H_2 + C^+ => H + CH^+$ (endothermicity: 4300 K)

 $H + CH^+ + => H_2 + C^+$

To overcome the energy threshold to form CH⁺, H₂ ro-vibrational energy can be used when available.

CH⁺ correlation with H₂^{*} suggests that the formation process of CH⁺ depends on H₂^{*} (in agreement with Lambert <mark>& Danks 1986,</mark> Agundez et. al 2010)

OH formation: $H_2 + O => H + OH$

OH comparison with H₂:

OH destruction: OH + photon => H + O $H + OH => H_2 + O$

The non-correlation between OH and H₂* suggests that the formation process of OH does not depend on H₂³ (in agreement with Agundez et. al 2010).



As well as corresponding to the irradiated structures (Goicoechea et al. 2011), the OH emission seems to also correspond with a bright young object, identified as a proplyd (see the zoomed image).

Physical properties of the Orion bar



<u>See also talk by E. Habart on</u> WED 16.10. at 12:20 (session 7b): **PET AL. Diagnostics as Observed** by PACS and SPIRE

We have selected a few positions across the bar and derive the physical properties using the radiative transfer code RADEX.

The best fits were found with H₂ density of Ix10⁵ and Ix10⁶

The temperature is higher closer to the ionizing stars (in front of the bar), and decreases further away from it. Column density is lower in front and behind the bar and higher in the bar.

The CO fit produces a high thermal pressure and small length, which confirm the presence of bright and dense structures in the bar.

The collisions can explain the excitation of warm CO, thus we conclude that the photoelectric heating is the dominant heating mechanism.



Ongoing work: We will use the constraints given by the RADEX models for input into PDR Meudon code. For the isobar model with P ~3×10⁸ K cm-3, the model published in Joblin et al. (in prep.) reproduces all the observed CO line intensities towards the Bar at the CO⁺ peak position.

REFERENCES: Agundez et al. (2010), ApJ, 713, 662 Bernard-Salas et al. (2012), A&A, 538, A37

Goicoechea et al. (2011), A&A, 530, L16 Joblin et al. (in prep.) Köhler et al. (2013, submitted)

Lambert & Danks (1986), ApJ, 303, 401 Lebouteiller et al. (2012), A&A 548, A91 Walmsley et al. (2000), A&A, 364, 301