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The Central 1000 au of a Pre-stellar Core Revealed with ALMA. II. Almost Complete Freeze-out

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ABSTRACT

Pre-stellar cores represent the initial conditions in the process of star and planet formation. Their low temperatures (<10 K) allow the formation of thick icy dust mantles, which will be partially preserved in the future protoplanetary disks, ultimately affecting the chemical composition of planetary systems. Previous observations have shown that carbon- and oxygen-bearing species, in particular CO, are heavily depleted in pre-stellar cores due to the efficient molecular freeze-out onto the surface of cold dust grains. However, N-bearing species such as NH₃ and, in particular, its deuterated isotopologues, appear to maintain high abundances where CO molecules are mainly in solid phase. Thanks to ALMA, we present here the first clear observational evidence of NH₂D freeze-out toward the L1544 pre-stellar core, suggestive of the presence of a “complete-depletion zone” within a $\simeq 1800$ au radius, in agreement with astrochemical pre-stellar core model predictions. Our state-of-the-art chemical model coupled with a non-LTE radiative transfer code demonstrates that NH₂D becomes mainly incorporated in icy mantles in the central 2000 au and starts freezing-out already at $\simeq 7000$ au. Radiative transfer effects within the pre-stellar core cause the NH₂D($1_{11} - 1_{01}$) emission to appear centrally concentrated, with a flattened distribution within the central $\simeq 3000$ au, unlike the 1.3 mm dust continuum emission which shows a clear peak within the central $\simeq 1800$ au. This prevented NH₂D freeze-out to be detected in previous observations, where the central 1000 au cannot be spatially resolved.

Keywords: Star formation (1569); Interstellar medium (847); Molecular clouds(1072)

1. INTRODUCTION

The initial conditions of the process of star and planet formation are to be found in dense cloud cores (Bergin & Tafalla 2007), especially in the so-called pre-stellar cores, which are gravitationally bound objects (André et al. 2014) with central number densities above 10^5 cm^{-3} (Keto & Caselli 2008) and large deuterium fractions

measured especially in N-bearing species (Crapsi et al. 2005), which preferentially trace the central regions (Hily-Blant et al. 2010; Spezzano et al. 2017). Detailed studies of pre-stellar cores in nearby clouds have provided clues on their physical and chemical structure: they are centrally concentrated, with low central temperatures (~ 7 K, Crapsi et al. 2007; Pagani et al. 2007; Launhardt et al. 2013) and evidence of grain growth (Forbrich et al. 2015; Chacón-Tanarro et al. 2019). They display a clear chemical differentiation, where rare CO isotopologue emission maps present a valley within the central ~ 8500 au, due to freeze-out onto dust grains (Caselli et al. 1999), while N-bearing species such as N_2H^+ , NH_3 and their deuterated forms appear to trace well the core center (Tafalla et al. 2002, 2004; Juárez et al. 2017; Spezzano et al. 2017), defined as the millimeter and sub-millimeter dust continuum emission peak. Although there is evidence of some central depletion of N_2H^+ (Bergin et al. 2002), N-bearing deuterated species have been found to trace well the dust peak (Caselli et al. 2002). Only recently, high-sensitivity single-dish observations of multiple transitions of N_2D^+ , coupled with astrochemical and non-LTE radiative transfer modeling, have provided evidence of N_2D^+ central depletion (Redaelli et al. 2019), but the size of the related depletion zone cannot be measured because of the limited angular resolution. The abundance of N_2H^+ isotopologues is linked to N_2 , which has a binding energy similar to that of CO (Bisschop et al. 2006), although recent studies have shown that the N_2 (but not the CO) binding energy is reduced in CO- N_2 ice mixtures (Nguyen et al. 2018). The smaller amount of N_2 freeze-out compared to CO can be explained if N_2 formation proceeds at a slower pace than CO formation, so that most of N_2 is still forming from (the more volatile) atomic nitrogen in the dense material when CO has already started to condense onto icy mantles (Flower et al. 2006; Hily-Blant et al. 2010; Pagani et al. 2012).

No evidence of ammonia freeze-out in starless and pre-stellar cores has been reported yet, despite the fact that the NH_3 binding energy is significantly larger than that of CO and N_2 and close to that of water molecules (Sipilä et al. 2019). This is puzzling, as the freeze-out time scale within the central 1000 au of pre-stellar cores, where the number density is about 10^6 cm^{-3} , is only ~ 1000 yr, much shorter than dynamical time scales. Furthermore, it appears that the NH_3 abundance is slightly *increasing* toward the center when viewed with single-dish and interferometers (Tafalla et al. 2002; Crapsi et al. 2007; Caselli et al. 2017). Such an increasing abundance profile for NH_3 is at odds with state-of-the-art astrochemical modeling (Sipilä et al. 2019). Crapsi et al. (2007)

also observed para- $\text{NH}_2\text{D}(1_{11} - 1_{01})$ toward the Taurus pre-stellar core L1544 with the IRAM Plateau de Bure Interferometer, finding an integrated intensity map centered on the dust peak, thus ruling out any central depletion of NH_2D when viewed at an angular resolution of $5.8'' \times 4.5''$ ($986 \text{ au} \times 765 \text{ au}$, at the distance of 170 pc¹; Galli et al. 2019). From this result, it was concluded that no *complete depletion zone*² (Walmsley et al. 2004) was present in L1544.

With ALMA 1.3 mm dust continuum emission observations of L1544, we have recently revealed a compact central region, with mass $\simeq 0.16 M_\odot$, radius $\simeq 1800$ au and average H_2 number density of $\simeq 10^6 \text{ cm}^{-3}$, called the “kernel” (Caselli et al. 2019). This kernel is the stellar system cradle, which we are investigating in detail to shed light on the first steps toward star and protoplanetary disk formation. Here we present the ALMA (12m+ACA) para- $\text{NH}_2\text{D}(1_{11} - 1_{01})$ mosaic, which reveals for the first time the NH_2D depletion zone toward the L1544 kernel at an angular resolution of $2.5''$ (425 au; Sections 2 and 3). We also demonstrate that the current data are in agreement with our astrochemical model predictions (Section 4), implying that almost all (99.99%) species heavier than He reside on dust icy mantles within the kernel. Hereafter, p NH_2D indicates para- NH_2D . A detailed study of the kernel kinematics will be presented in a forthcoming paper.

2. OBSERVATIONS

2.1. Deuterated Ammonia

We observed L1544 with ALMA during Cycle 4 (ESO Project ID 2016.1.00240.S, PI Caselli) using the 12 m (Main array) and the 7 m (ACA) arrays. The 12 m array 3-pointing observations were carried out on 2016 December 22, while the 7 m array single pointing observations were carried out on 2016 October 2. The Main array observations included the quasar J0510+1800 for bandpass and phase calibration, while quasar J0423-0120 is observed for flux calibration. The ACA observations included the quasar J0510+1800 for flux, bandpass, and phase calibration. The 3-point mosaic observations used a correlator with the continuum window centered at 100 GHz and a spectral window for p $\text{NH}_2\text{D}(1_{11}-1_{01})$ centered at 110.15355 GHz (these data will be presented

¹ We note that previous papers on L1544 adopted a distance to the source of 140 pc or 135 pc, based on previous distance measurements (Schlafly et al. 2014; Galli et al. 2018). Here we adjust previous results to the new distance estimate.

² A *complete depletion zone* is defined as a region within a pre-stellar core where, due to freeze-out, species heavier than He have left the gas phase and reside in the icy mantles of dust grains.

in a future paper, dedicated to the study of dust properties within the kernel). The bandwidth and spectral resolution for the pNH_2D observations are 58.59 MHz and 30.518 kHz (corresponding to a velocity resolution of 83 m s^{-1}), respectively. The 12 m and 7 m array data were calibrated using the Common Astronomy Software Applications package (CASA McMullin et al. 2007) version 4.7.0. We assigned weights to the datasets with the `statwt` command before combining them (`concat` command). We perform a joint deconvolution of both data sets (ACA and 12 m array) in CASA (version 5.6.2) using `tclean` for mosaics, natural weighting, and restored with a circular beam of $2.5''$. We use a rest frequency for the imaging of 110.153562 GHz (Cohen & Pickett 1982; De Lucia & Helminger 1975). We use the multi-scale clean technique, with the multi-scale parameter of $[0, 2.5, 7.5, 22.5]''$. The final noise level is 7 mJy beam^{-1} per 0.08 km s^{-1} channel, which corresponds to a noise level of 0.13 K in the central $45''$ diameter region of the primary beam corrected data. The integrated intensity map of $\text{pNH}_2\text{D}(1_{11} - 1_{01})$ (ACA+12 m array) is shown in the top panel of Fig. 1 and the full mosaic is shown in Fig. A1. We note that the combined ACA+12 m data perfectly recover the single dish flux within the IRAM-30m beam (see Figure B1), demonstrating that mosaic pointing observations with ALMA are a powerful technique to carry out detailed studies of pre-stellar cores.

2.2. 1.3 mm Continuum

We observed L1544 with ALMA during Cycle 2 (ESO Project ID 2013.1.01195.S, PI Caselli) using both the 12 m (Main array) and the 7 m (ACA) arrays. The data were already presented in Caselli et al. (2019), but here we briefly describe the data reduction process. The 12 m array observations were carried out on 2014 December 27 and 29, while the 7 m array observations were carried out on 2014 June 15, July 20 and 29, and August 6 and 11. The single pointing observations used a correlator with the continuum centered at 228.973 GHz and a 2 GHz bandwidth. The 12 m array data were calibrated using the CASA version 4.2.2, while the 7 m array data were calibrated using CASA version 4.2.1. We run the `statwt` command on both datasets before combination. We perform a joint deconvolution of both data sets (ACA and 12 m array) in CASA (version 5.6.2) using `tclean` multifrequency synthesis (mfs) for mosaics, a natural weighting, a taper of $2.1'' \times 1.45''$ and PA of 147.2° , and restored with a circular beam of $2.5''$. Since the continuum emission is extended, we use the multi-scale clean technique, with the multi-scale parameter of $[0, 1.25, 3.75, 11.25]''$. The 1.3 mm dust continuum emission map can be seen in the bottom panel of Fig. 1,

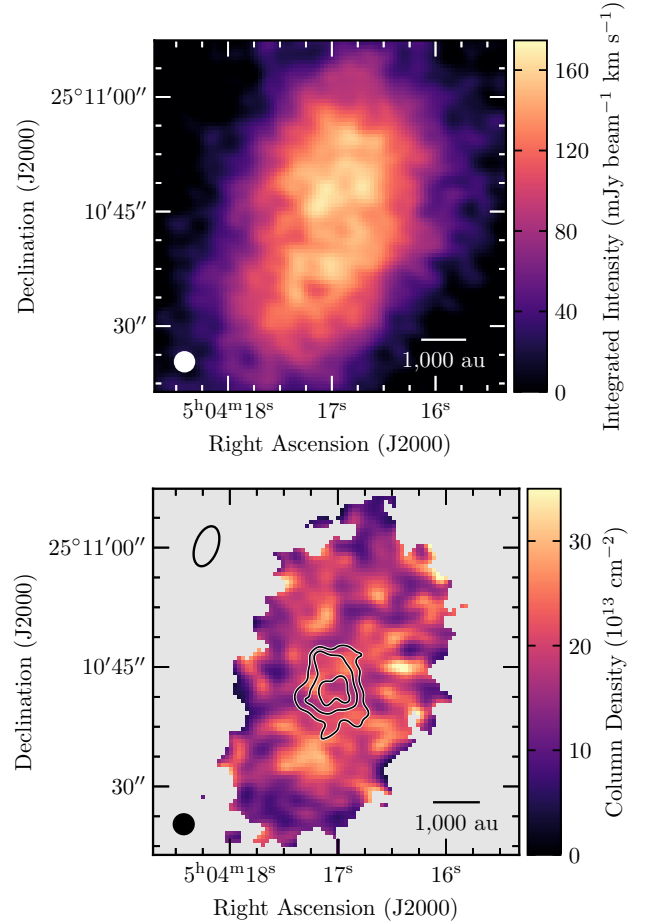


Figure 1. *Top:* Integrated intensity map of the $\text{pNH}_2\text{D}(1_{11} - 1_{01})$ transition toward L1544. *Bottom:* pNH_2D column density map derived from fitting the hyperfine structure. The contours show the 14, 17, and $20 \times \sigma$ levels ($\sigma = 36 \mu\text{Jy beam}^{-1}$) of the 1.3 mm dust continuum emission map presented in Caselli et al. (2019), which locates the kernel. Scale bar and beam size are shown at the bottom right and left corners, respectively. In the top left of the bottom panel we show the orientation and shape of the ellipse used to define the projected radius in Sect. 3 and Fig. 2. The typical uncertainty on the pNH_2D column density is about 10%.

overlaid onto the pNH_2D column density map, which will be presented in Sect. 3.

3. COLUMN DENSITY CALCULATION

We perform the line fit of $\text{pNH}_2\text{D}(1_{11} - 1_{01})$ using the `nh2d` model implemented in `pyspeckit` (Ginsburg & Mirocha 2011). The $\text{pNH}_2\text{D}(1_{11} - 1_{01})$ line is modeled using the hyperfine structure and relative intensity from the recent laboratory works of Melosso et al. (2020) and Melosso et al. (2021). The model assumes a Gaussian velocity distribution (with centroid velocity and velocity dispersion V_{lsr} and σ_v), equal excitation temperatures (T_{ex}) for all hyperfine components, and total op-

tical depth (τ_0). We perform a fit for the whole cube with all 4 free parameters. Following the procedure in Pineda et al. (2021), we discard all centroid velocity determinations if the uncertainty in the V_{lsr} is larger than 0.02 km s^{-1} , the velocity dispersion determinations if the uncertainty σ_v is larger than 0.015 km s^{-1} , and the excitation temperature if the uncertainty is larger than 1 K.

The column density has been derived following Mangum & Shirley (2015):

$$N(\text{pNH}_2\text{D}) = \frac{8\pi\nu^3}{c^3 A_{ul}} \frac{Q(T_{ex})}{g_u e^{-E_u/kT_{ex}}} \times \left[e^{h\nu/kT_{ex}} - 1 \right]^{-1} \int \tau dv, \quad (1)$$

where $\int \tau dv = \sqrt{2\pi} \tau_0 \sigma_v$, ν is the frequency of the transition, g_u is the upper level degeneracy, E_u is the upper energy level of the transition, A_{ul} is the Einstein coefficient, and $Q(T_{ex})$ is the partition function. The values for the E_u , g_u , and A_{ul} are obtained from the Leiden Atomic and Molecular Database (LAMDA; Schöier et al. 2005; van der Tak et al. 2020)³, and the partition function is calculated up to the first 30 energy levels. This formula has been used previously (Daniel et al. 2016; Harju et al. 2017). The T_{ex} has been estimated at each position using the hfs fitting within `pyspeckit`, with average uncertainties of 2% (and a maximum of 5% at the edge of the map). The error associated with the column density is about 10%, close to the calibration error.

The column density map in Fig. 1 shows a complex structure, partly due to noise, without a clear peak toward the 1.3 mm dust continuum peak, which identifies the kernel (Caselli et al. 2019). This implies that the pNH₂D column density does not follow the H₂ column density, thus a drop in pNH₂D fractional abundance (with respect to H₂) within the kernel radius of $\simeq 1800 \text{ au}$ ($\simeq 10''$) must be present. The pNH₂D column density flattening within the central $\simeq 3000 \text{ au}$, which includes the L1544 kernel, is well illustrated in Fig. 2. The circles represent individual measurements, while the black and gray horizontal lines are the mean and its associated uncertainty in a $2.5''$ bin. The projected radius pr is calculated as $pr = \sqrt{r_{maj}^2 + (r_{min} b_{axes})^2}$, where r_{maj} and r_{min} are the distances along the semi-major and -minor axes, respectively; $b_{axes} = 22/12$ is the semi-major and -minor axis ratio for the pNH₂D emission, where the semi-major axis is along the position angle of 160° (measured East from North; see the bottom panel

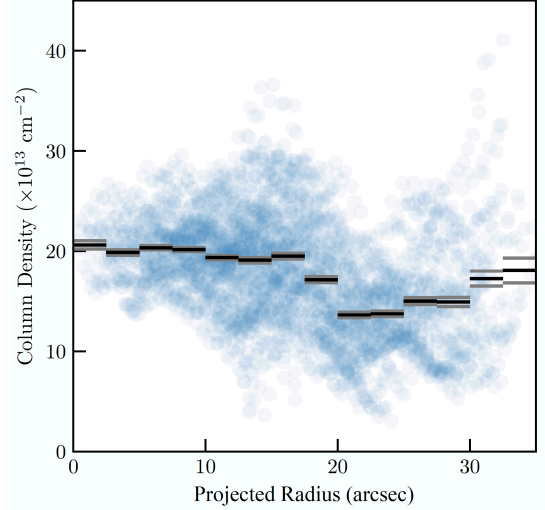


Figure 2. pNH₂D column density as a function of projected radius (see main text for its definition). Blue circles correspond to individual measurements across the column density map in the bottom panel of Fig. 1. The black and gray horizontal lines are the mean and corresponding uncertainty within a $2.5''$ bin. Note the flat column density profile within the central $10''$, where the 1.3 mm dust continuum map shows a clear peak (see Fig. 1, bottom panel).

of Fig. 1 for an illustration of the elliptical shape and orientation used to define pr).

4. CHEMICAL AND RADIATIVE TRANSFER MODELING

We carried out chemical and radiative transfer modeling to reproduce the observed pNH₂D column density and line emission distribution. First, we produced simulated pNH₂D abundance distributions using the most recent version of our gas-grain chemical code (Sipilä et al. 2020), which includes an extensive description of deuterium and spin-state chemistry, of critical importance for the present work. Very briefly, we used the one-dimensional (1D) source model for L1544 presented by Keto et al. (2015), which was divided into concentric shells; chemical simulations were carried out separately in each shell. This process yields time-dependent abundance profiles. The initial abundances used in the chemical modeling are displayed in Table 1. We assume a constant grain radius of $0.1 \mu\text{m}$, and a grain material density of 2.5 g cm^{-3} . Other model details can be found in Sipilä et al. (2020) and are omitted here for brevity.

We searched for the best fit to the observed line map by extracting the pNH₂D abundance profile from the chemical model at various time steps (between 10^4 and 10^6 yr), and using the abundance data as input for the non-LTE line radiative transfer code LOC (Juvela 2020) to simulate the line emission. We adopted the pNH₂D –

³ <https://home.strw.leidenuniv.nl/~moldata/datafiles/p-nh2d.dat>

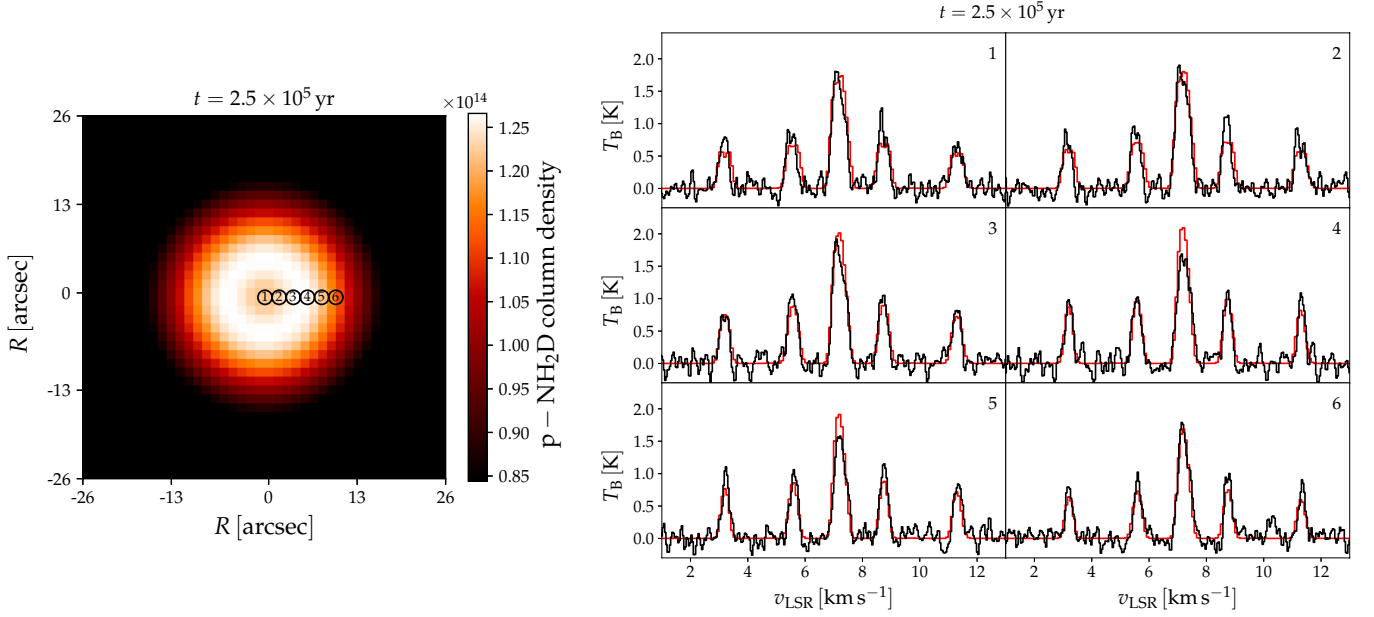


Figure 3. *Left:* Map of the pNH₂D column density distribution predicted by the chemical model at the best-fit time. The column densities have been convolved to a beam of 2.5'' (425 au). The black circles with numbers are the areas within which the red spectra in the right panel, with the corresponding numbers, have been extracted. *Right:* Simulated pNH₂D (1₁₁ – 1₀₁) line emission profiles (red) at $\sim 2.5''$ intervals from the core center. The observed lines toward equivalent positions in the South-East direction in Figure A1, parallel to the major axis, are shown in black.

Table 1. Initial fractional abundances (with respect to total H nuclei, n_{H}).

Species	Abundance
H ₂	5.00×10^{-1} ^(a)
He	9.00×10^{-2}
C ⁺	1.20×10^{-4}
N	7.60×10^{-5}
O	2.56×10^{-4}
S ⁺	8.00×10^{-8}
Si ⁺	8.00×10^{-9}
Na ⁺	2.00×10^{-9}
Mg ⁺	7.00×10^{-9}
Fe ⁺	3.00×10^{-9}
P ⁺	2.00×10^{-10}
Cl ⁺	1.00×10^{-9}

^aThe initial H₂ ortho/para ratio is 1×10^{-3} .

pH₂ collisional coefficients from Daniel et al. (2014) and assumed that the hyperfine components in each rotational level have the same excitation temperature. The hyperfine component frequencies and relative intensities have been adopted from Melosso et al. (2021). The velocity profile of the contracting cloud (including a uniform level of turbulence of 0.1 km s^{-1}) was taken from Keto et al. (2015). With this setup, the best fit between the model and the observations is reached after a chem-

ical evolution of $t \sim 2.5 \times 10^5 \text{ yr}$, as determined by a χ^2 analysis applied to modeled and observed spectra toward six positions at and near the center of the core⁴. We stress that the chemical best-fit time of $2.5 \times 10^5 \text{ yr}$ cannot be used as a chemical clock or to deduce the dynamical age of the core. The reason for this is that the chemistry is run using a static physical structure as input (a BE sphere with a density, temperature and velocity profile deduced by Keto et al. 2015), so chemical processes are not followed simultaneously with the dynamical evolution. Future work will focus on self-consistent chemical-dynamical evolution models, using hydro- and magneto-hydrodynamics simulations, of the types mentioned in Appendix D.

Figure 3 shows the simulated pNH₂D column density map and the (1₁₁ – 1₀₁) line emission profiles at the best-fit time. The modeled column densities and line emission maps agree well with the observations, showing a relatively flat column density map within the central

⁴ The six observed spectra have been extracted from the six 2.5'' circular areas, along the major axis and toward the South-East direction, displayed in Fig. A1. We also compared with the six spectra extracted from the perpendicular direction along the minor axis (see Fig. A1) and found similar results within the central 1800 au of the kernel (see Appendix A and Fig. A2), while the outer positions show some deviation, suggesting that the (spherically symmetric) model is in better agreement with the physical structure along the major axis (see also Appendix D).

$\simeq 13''$ (or about 2200 au). In fact, although the pNH_2D column density map in Fig. 3, left panel, shows a central valley within a bright ring of $\simeq 5''$ ($\simeq 850$ au), while our observations show a flattened structure (see Fig. 2), the model column density at the dust peak is only 4% lower than at the bright ring ($1.22 \times 10^{14} \text{ cm}^{-2}$ instead of $1.27 \times 10^{14} \text{ cm}^{-2}$), well within our calibration uncertainties. As we will show in Section 5, this implies that our observations are consistent with NH_2D freeze-out in the central core. We have verified that the $(1_{11} - 1_{01})$ line emission profiles cannot be reproduced by the models without NH_2D depletion toward the center (see Appendix C). We note that the observed average pNH_2D column density in Fig. 2 is slightly (less than a factor of two) larger than that in Fig. 3 ($\simeq 2 \times 10^{14} \text{ cm}^{-2}$ versus $\simeq 1.2 \times 10^{14} \text{ cm}^{-2}$), which is once again underlying the good agreement between model and observations, despite the simplicity of the model. The measured and modeled excitation temperatures T_{ex} are also similar, as can be seen in Fig. 4, which compares the measured T_{ex} (as deduced from the `pyspeckit` hfs fit of all the spectra within the map in Fig. 1) as a function of *projected radius* (pr , see Sect. 3) with the excitation temperature profile *within* the core ($T_{\text{ex}}(r)$, with r the core *radius*) predicted by the radiative transfer LOC. Although a one-to-one comparison between the two excitation temperatures cannot be made, as one should instead compare the measured T_{ex} with the integral along the line of sight of the model $T_{\text{ex}}(b)$ for each impact parameter b , the two values are comparable, with the model $T_{\text{ex}}(r)$ within 20-30% the measured T_{ex} .

A final note concerns the adopted physical structure. The Keto et al. (2015) 1D model used here provides self-consistently volume density and velocity profiles across the core and it can be easily input in comprehensive gas-grain chemical models such as the present one and previous work by, e.g., Vasyunin et al. (2017) on complex organic molecules in L1544. Naturally, this 1D model cannot reproduce the elongated structure observed at 1.3mm (see Figure 3 in Caselli et al. 2019, and Appendix D). In fact, ALMA observations are better reproduced by 3D non-ideal magneto-hydrodynamic (niMHD) simulations of a contracting cloud with peak volume density of 10^7 cm^{-3} (see their Figure 4). The 3D niMHD simulation gives rise to a pre-stellar core with an elongated flattened central structure, in harmony with the observed 1.3mm ALMA dust continuum emission. The adoption of this 3D model for the chemistry is however beyond the scope of this paper and it will be used in a future publication, where only simple species will be followed in a reduced chemical network to be included in the dynamical simulation. Despite the

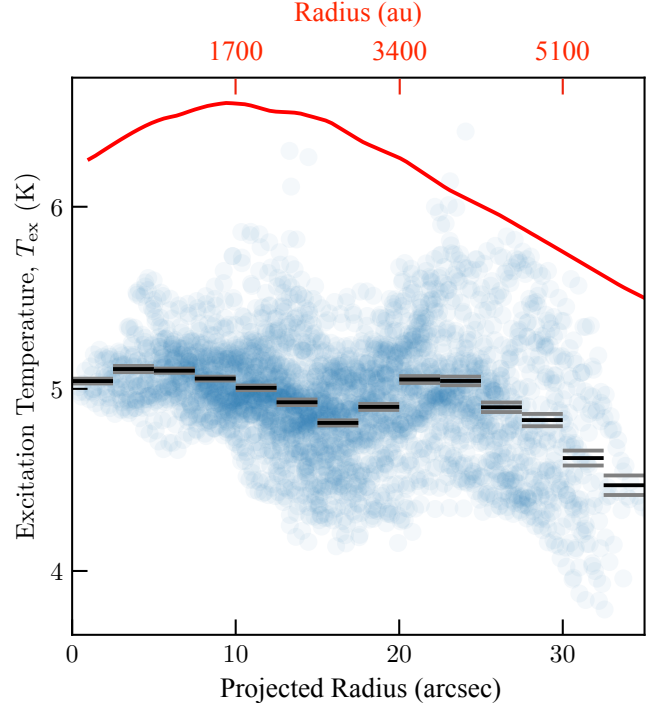


Figure 4. $\text{pNH}_2\text{D}(1_{11} - 1_{01})$ excitation temperature (T_{ex}) as a function of *projected radius* (see Section 3 for its definition). Blue circles correspond to individual measurements. The black and gray horizontal lines are the mean and corresponding uncertainty within a $2.5''$ bin. The red curve is the model T_{ex} profile, the predicted excitation temperature as a function of core *radius*, $T_{\text{ex}}(r)$.

difference of the two (1D and 3D) models, it is interesting to note that the density and velocity profiles along the major and minor axis of the simulated 3D core do not differ substantially from those predicted by the 1D model of Keto et al. (2015), as shown in Fig. D1. In particular, the density and velocity profiles along the major axis of the 3D model are especially close to our adopted 1D model (see Fig. D1), suggesting similarities that, together with the radiative transfer results shown in Fig. 3, make our 1D model a very good approximation of the L1544 structure.

5. DISCUSSION: ALMOST COMPLETE FREEZE-OUT IN THE KERNEL

In the previous sections we showed that ALMA has allowed us to resolve for the first time the distribution of deuterated ammonia in the central 1000 au of the pre-stellar core L1544. Despite the complex morphology of the L1544 core as viewed at the frequency of the $\text{pNH}_2\text{D}(1_{11} - 1_{01})$ line (see Fig. 1), it is clear that the pNH_2D column density has a flattened distribution in correspondence of the dust continuum peak, where the kernel is located (see also Fig. 2). The radius of the kernel is

about $\simeq 1800$ au, which is ~ 5 times smaller than the CO depletion zone observed toward the same pre-stellar core by Caselli et al. (1999). This is the reason why previous interferometric observations from Crapsi et al. (2007), done with poorer angular resolution ($5.8'' \times 4.5''$), could not reveal the abundance drop of deuterated ammonia, as demonstrated in Appendix E.

In the following, we will present the fractional abundance profiles predicted by our chemical model, at 2.5×10^5 yr, of all ammonia isotopologues as well as the profile of the *total depletion factor*, $f_{D_{\text{tot}}}$, defined by:

$$f_{D_{\text{tot}}} = \frac{\sum_i [x(i)_{\text{gas}} + x(i)_{\text{solid}}]}{\sum_i x(i)_{\text{gas}}}, \quad (2)$$

where i represents any species in the chemical model containing an element heavier than He. Thus, $f_{D_{\text{tot}}}$ quantifies the amount of elements heavier than He subsisting in the gas phase, with the higher values corresponding to higher levels of depletion.

The abundance profile of pNH_2D at the best-fit time is displayed in Fig. 5 along with $f_{D_{\text{tot}}}$. It is interesting to note that the pNH_2D abundance starts to drop already within about 7000 au, whereas the integrated intensity map as well as the simulated observations show a centrally concentrated structure up to the central $\simeq 2000$ au. The abundance drop is in fact partially hidden to observations due to the radiative transfer of the $\text{pNH}_2\text{D}(1_{11} - 1_{01})$ line. In particular, despite a ‘‘catastrophic’’ pNH_2D freeze-out present within the central 2000 au (see Fig. 5), the corresponding decrease in column density is only $\sim 4\%$; this is shown in the left panel of Fig. 3, where a relatively bright ring of about $5''$ in radius, within the pNH_2D depletion zone, is clearly visible. This is due to the fact that the column density is dominated by larger scales, where pNH_2D is abundant. The critical density of the $(1_{11} - 1_{01})$ line is $\sim 1.4 \times 10^5 \text{ cm}^{-3}$ at $T = 10$ K, and so collisional excitation dominates in a region roughly $13''$ in radius, enclosing the bright ring in the left panel of Fig. 3. The inferred column density ring and shallow inner valley (Fig. 3) are due to the drop in the pNH_2D abundance toward the center of the core. The simulated main hyperfine component peak has an optical depth close to 1, in agreement with our observations, and the strength of the whole line emission follows the morphology of the column density distribution.

The abundance distributions of the modeled normal and deuterated ammonia (summed over the spin states) as well as that of CO are also shown in Fig. 5 (right panel); it is evident that the depletion of singly-deuterated ammonia within the central 2000 au is not due to conversion into multiply-deuterated ammonia, but is instead a result of freeze-out onto grain sur-

faces. These results, including the ortho-to-para ratios of NH_2D , are consistent with those obtained by Hily-Blant et al. (2018), who modelled the molecular composition and the nuclear spin chemistry of a collapsing Bonnor-Ebert (BE; Bonnor 1956; Ebert 1955) sphere. According to our model, the total depletion factor $f_{D_{\text{tot}}}$ increases from 1000 to 10000 in the central 2000 au, implying that between 99.9% and 99.99% of all the species heavier than He are predicted to reside on the surface of dust grains within the kernel. This has the consequence that thick icy mantles are expected to form on top of dust grains within the central regions of pre-stellar cores, just before the formation of a protostar and protostellar disk. Thick icy mantles could be responsible for the change in dust opacity measured toward the center of L1544 by Chac3n-Tanarro et al. (2019).

We note that, previous to this work, coarser interferometric observations of pNH_2D showed a clear centrally peaked distribution toward L1544 (Crapsi et al. 2007). For this reason, pNH_2D was considered the species most resilient to freeze-out (compared with all other species observed so far toward this core which all show direct or indirect evidence of depletion, including N_2D^+ ; e.g. Spezzano et al. 2017; Redaelli et al. 2019). Now that even pNH_2D freeze-out has been unveiled and our modeling reproduces the observed column density map and line profiles of $\text{pNH}_2\text{D}(1_{11} - 1_{01})$, we can conclude that the almost complete freeze-out predicted by our chemical model is then consistent with the data available for this source. Moreover, as already mentioned, previous single-dish and interferometric observations suggested a NH_3 abundance profile increasing toward the central regions (e.g., Crapsi et al. 2007). To investigate this apparently contradicting result, we simulate also the $\text{NH}_3(1,1)$ line profiles using the pNH_3 abundance profile predicted by our model, after 2.5×10^5 yr of chemical evolution, and compare them with the VLA-only observations in Figure 2 of Crapsi et al. (2007). We also consider the case without NH_3 freeze-out in the central 6000 au (which differs from our predicted profile shown in Fig. 5, right panel) and see if this better reproduces the observations, in line with the conclusions from Crapsi et al. (2007). The assumed beam for the simulated observations of $\text{NH}_3(1,1)$ is $4''$, close to the VLA resolution of $4.35'' \times 3.45''$. The spectra are reported in Fig. F1, showing that our model prediction of NH_3 freeze-out toward the center (see Fig. 5) is consistent with the Crapsi et al. (2007) observations, including the relative intensity of the groups of hyperfines and the negligible difference between the spectrum centered at the dust peak position and the one at the $22.3''$ offset. The main difference is that our simulated lines are

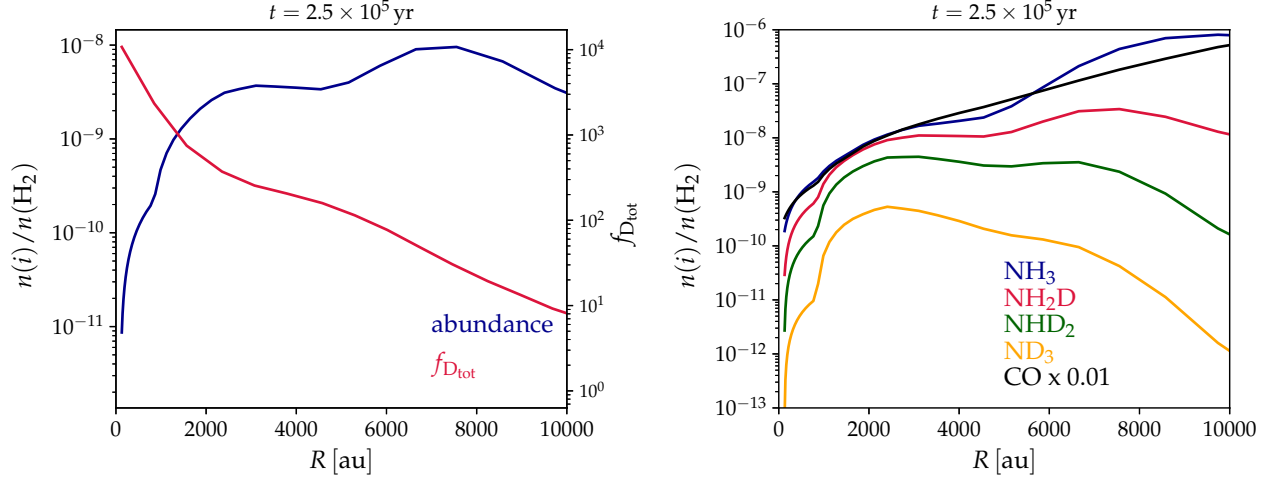


Figure 5. *Left:* Model fractional abundance of pNH_2D (blue) and overall gas-phase depletion factor (red) at the best-fit time as a function of radius. *Right:* Fractional abundances w.r.t. H_2 of normal and deuterated ammonia (colors, indicated in the Figure), as well as of CO (multiplied by 0.01, black), at the best-fit time as a function of distance from the core center. The ammonia fractional abundances represent sums over the nuclear spin forms (ortho and para; also meta for ND_3).

about two times brighter than the observed ones, but this is expected due to the fact that we are not simulating interferometric observations, and the core envelope is significantly contributing to the simulated line flux (unlike in the case of interferometric observations). In fact, higher sensitivity VLA data of $\text{NH}_3(1,1)$ toward L1544, combined with GBT single-dish data, show a peak temperature in better agreement with our modeling; these data will be presented in an upcoming paper focusing on a new gas temperature map of L1544 (Schmiedeke et al., in prep.). Another interesting result of this comparison is that variations in the abundance profile of ammonia in the central 5000 au do not change the spectra (see dashed histograms in Fig. F1, which refer to a model with a flat fractional abundance of NH_3 of 2×10^{-8} within the central 5000 au), demonstrating that $\text{NH}_3(1,1)$ cannot be used to measure its abundance profile within the central few thousand au, because of its relatively low critical density (see Appendix F). Interestingly, new high-resolution and high-sensitivity NH_3 observations of the pre-stellar core H-MM1 in Ophiuchus show a clear depletion zone towards its dust peak (Pineda et al., in prep.), thus confirming our predictions.

A final comment should be added about the velocity structure used as input to the chemical model. In some previous studies of L1544, the Keto & Caselli (2010) infall velocity profile had to be scaled up by a factor of ~ 2 to reproduce observations obtained with single-dish telescopes (Bizzocchi et al. 2013; Redaelli et al. 2019). In the present case, a similar scaling leads to double-peaked line profiles, even for the satellite components, that are completely excluded by the observations. This discrep-

ancy with the earlier single-dish studies is likely related to the 1D spherically symmetric structure of the source model (while the source itself is certainly not spherical on the large scales traced by single-dish observations), and to the possible presence of more complex velocity fields at the core scale (see Appendix D). The present interferometric observations probe a relatively small volume, where the number density is close to the critical density of the observed transition ($1.4 \times 10^5 \text{ cm}^{-3}$), i.e. a region with radius of about 4000 au, considering the adopted density profile shown in Fig. D1 (red-dashed curve); hence the 1D symmetry assumption is adequate, as already discussed. Therefore, the Keto & Caselli (2010) physical model, which already successfully reproduced line profiles of several species observed with single-dish telescopes toward L1544 (e.g., Caselli et al. 2012; Keto et al. 2015), still remains valid for these new ALMA interferometric observations, as well as recent work from Redaelli et al. (2021).

6. CONCLUSIONS

Our ALMA spectroscopic observations, coupled with chemical and radiative transfer modeling, of the prototypical pre-stellar core L1544 have unveiled the depletion zone of deuterated ammonia, with size very close to the kernel previously found with ALMA continuum observations (radius $\simeq 1800$ au). This has not been possible before because of the poorer angular resolution (factor of ≥ 2) and quality of previous observations of the same line. Our chemical model, applied to the physical structure of L1544, reproduces remarkably well the $\text{pNH}_2\text{D}(1_{11} - 1_{01})$ emission. Simulated observations of

the modelled pre-stellar core show that the high pNH_2D fractional abundance in regions outside the kernel, with volume densities close to the critical density of the (1_{11} - 1_{01}) transition, prevent a clear view of the actual catastrophic NH_2D freeze-out region without high enough angular resolution. We find that more than 99.9% of all species heavier than He reside in thick icy mantles within the central kernel, the future stellar cradle. In pre-stellar cores like L1544, with central volume densities $\geq 10^6 \text{ cm}^{-3}$ and corresponding freeze-out time scales $\leq 10^3 \text{ yr}$, an almost complete freeze-out is therefore inevitable, just before star formation. The thick icy mantles (about 150 monolayers on the $0.1 \mu\text{m}$ -grains in our chemical model, close to the predictions from [Draine 1985](#)) may promote coagulation of dust grains (e.g., [Ormel et al. 2009](#); [Chacón-Tanarro et al. 2019](#)) and allow preservation and delivery of the frozen pre-stellar chemistry into the next stages of evolution, when the protostar and protoplanetary disk will form. Some of this pre-stellar ice, especially that trapped within icy pebbles in the outer part of the disk, may survive later stages of planet formation and evolution. Indeed, similarities in molecular composition and isotopic fractionation between star- and planet-forming regions and primitive material in our Solar System, such as comets and carbonaceous chondrites (e.g. [Mumma & Charnley 2011](#); [Caselli & Ceccarelli 2012](#); [Ceccarelli et al.](#)

[2014](#); [van Dishoeck et al. 2014](#); [Öberg et al. 2015](#); [Altwegg et al. 2019](#); [Cleeves et al. 2014](#); [Drozdovskaya et al. 2021](#)), strongly hint at a crucial role of pre-stellar chemistry in shaping the composition and further chemical/mineralogy evolution of the building blocks of planets.

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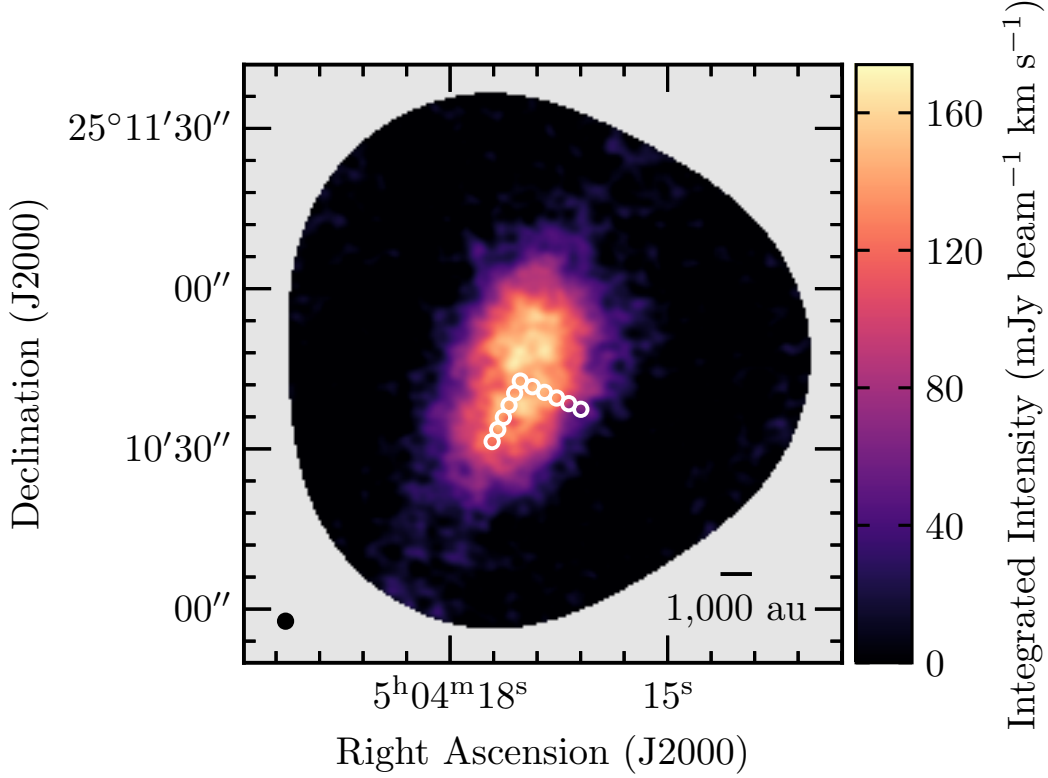


Figure A1. Integrated intensity map of the pNH₂D (1₁₁-1₀₁) transition on L1544, showing the 3-point mosaic. The white circles overlaid to the integrated intensity map mark the positions where the spectra of Fig. 3 (South-East direction, along the major axis) and Fig. A2 (South-West direction, along the minor axis) have been extracted.

APPENDIX

A. FULL MOSAIC MAP

In Figure 1, a zoom-in view of the pNH₂D(1₁₁ – 1₀₁) integrated intensity map is shown, in order to highlight the structure and better compare with the column density map. In this Appendix, we present the full 3-point mosaic of the same map to show the extent of the mapped area (see Fig. A1). The six white circles along the major axis in the South-East direction in Fig. A1 enclose the areas where the spectra have been extracted for comparison with the model spectra (see Fig. 3). The spectra from the other six white circles, those along the minor axis, are reported in Fig. A2 and compared with the same model spectra as in Fig. 3. It is apparent that the outer positions in Fig. A2 (numbers 5 and 6) show model spectra brighter than the observed ones, suggesting that the major axis has a physical and chemical structure closer to our spherically symmetric model. This is supported also by Fig. D1, where the density and velocity profiles of the 3D model are compared with our 1D model profiles.

B. COMPARISON WITH SINGLE DISH NH₂D(1₁₁ – 1₀₁) SPECTRUM

Figure B1 shows the comparison between the pNH₂D(1₁₁ – 1₀₁) spectrum obtained with the IRAM-30m single dish observations (with 22'' beam) and the one obtained by ALMA (12m+ACA) within the same area. The IRAM 30m spectrum of pNH₂D(1₁₁ – 1₀₁) has been extracted at the dust peak of an On-The-Fly map covering the inner 22'' × 22'' of L1544 around the dust peak. The emission map was obtained in the framework of the project 013-13 (PI: S. Spezzano) in October 2013. Position switching was used, with the reference position set at (-180'', 180'') offset with respect to the map centre. The EMIR E090 receiver was used with the Fourier Transform Spectrometer backend (FTS) with a spectral resolution of 50 kHz. The mapping was carried out in good weather conditions ($\tau \sim 0.03$) and a

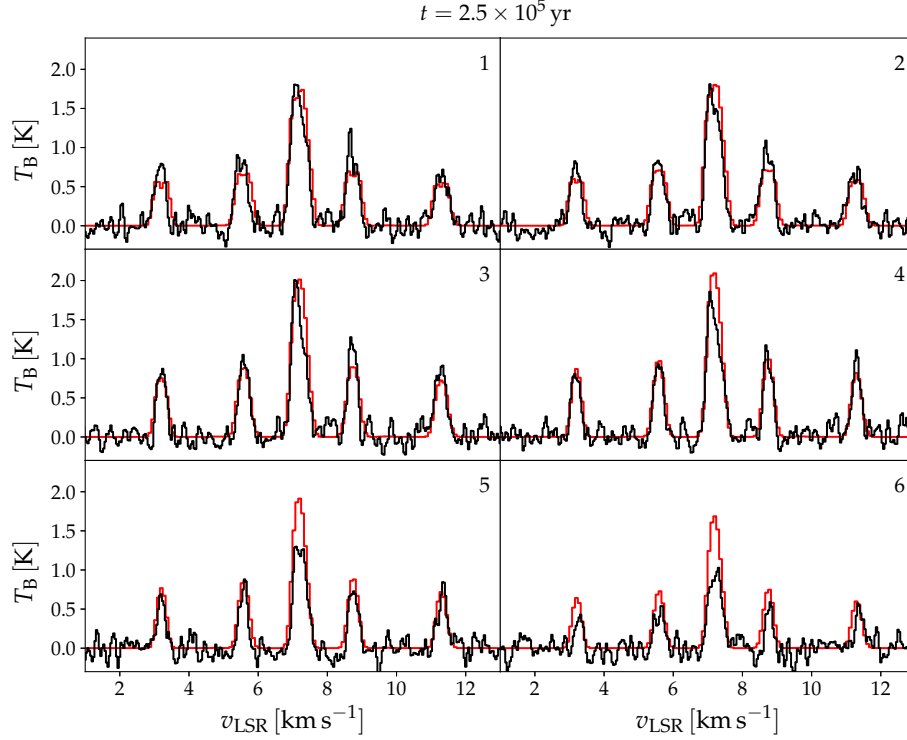


Figure A2. Same as the right panel of Fig. 3 but now the black histograms are the spectra extracted from the six circles in the South-West direction in Fig. A1 (along the L1544 minor axis). Note that the observed spectra in the outer positions (numbers 5 and 6) are now less intense than those predicted by our model. This suggests that the major axis of L1544 has physical and chemical structure in better agreement with our 1D model.

typical system temperature of ~ 90 - 100 K. It is clear from the figure that all the flux is fully recovered, demonstrating that ALMA mosaicing, together with the combination of 12 m and ACA data, provides a powerful technique to study pre-stellar cores in detail.

C. MODELED SPECTRA IN CASE OF NO CENTRAL FREEZE-OUT

As shown in Fig. 3, our model reproduces very well the $\text{pNH}_2\text{D}(1_{11} - 1_{01})$ spectra toward the L1544 kernel. This model predicts that NH_2D (as well as other species containing elements heavier than He) quickly disappears from the gas phase within the central 2000 au because of freeze-out onto dust grains (see Fig. 5). To test that central freeze-out is indeed needed to reproduce observations, Figure C1 shows the predicted $\text{pNH}_2\text{D}(1_{11} - 1_{01})$ spectra when the NH_2D fractional abundance is kept at the constant value of 6×10^{-9} within the central 2000 au, instead of dropping as in Fig. 5. The predicted spectra clearly overestimate the observed intensities in the central four positions, thus demonstrating that observations can only be reproduced when freeze-out is taken into account, as predicted by our chemical models.

D. DENSITY AND VELOCITY PROFILES

The physical structure of the L1544 core used for the chemistry code has been described in Sect. 4. As discussed there, the density profile from the self-consistent hydro-dynamic simulation of a contracting Bonnor-Ebert (BE; Bonnor 1956; Ebert 1955) sphere in quasi-static equilibrium (Keto & Caselli 2010) well reproduces line profiles previously observed toward the L1544 dust peak (e.g. Keto & Caselli 2010; Caselli et al. 2012; Keto et al. 2015). However, because of the spherical symmetry of the BE model, the elongated 1.3 mm dust continuum emission observed at high angular resolution with ALMA cannot be reproduced (Caselli et al. 2019). Indeed, Caselli et al. (2019) showed that the central kernel is consistent with flattening produced by non-ideal magneto-hydrodynamic (niMHD) simulations of the contraction of a rotating magnetized cloud core. The peak density within the flattened structure in the 3D niMHD simulation reaches 10^7 cm^{-3} , in agreement with the peak density at the center of the BE sphere, while the average density within the kernel is about 10^6 cm^{-3} , consistent with previous single-dish observations (Crapsi et al. 2007; Chacón-Tanarro et al. 2019), which did not have high enough angular resolution to detect the central density enhancement.

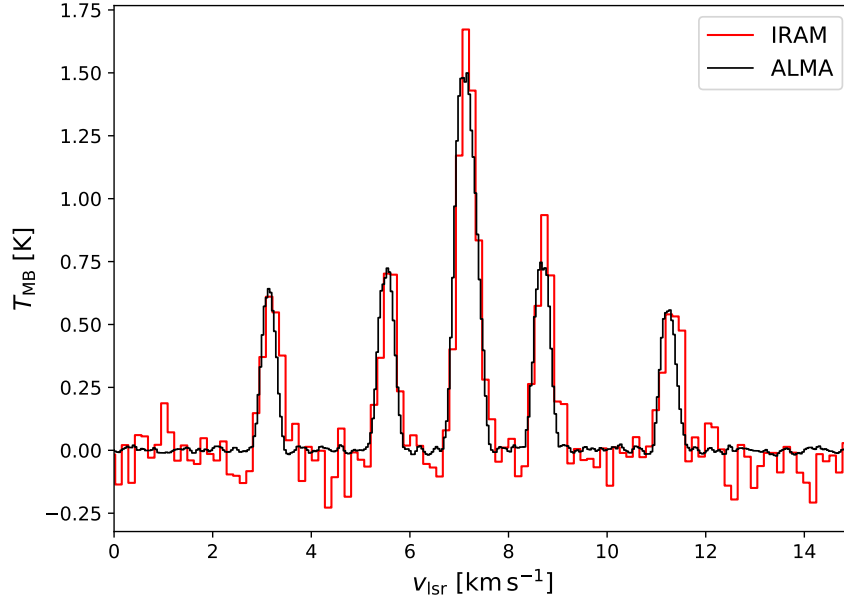


Figure B1. Comparison between the $\text{pNH}_2\text{D}(1_{11} - 1_{01})$ IRAM-30m spectrum (red histogram) and the ALMA 12m+ACA spectrum within the same $22''$ beam of the 30m observations (black histogram), obtained toward the L1544 kernel. This clearly shows that the ALMA 12m+ACA combined mosaic fully recovers the line flux.

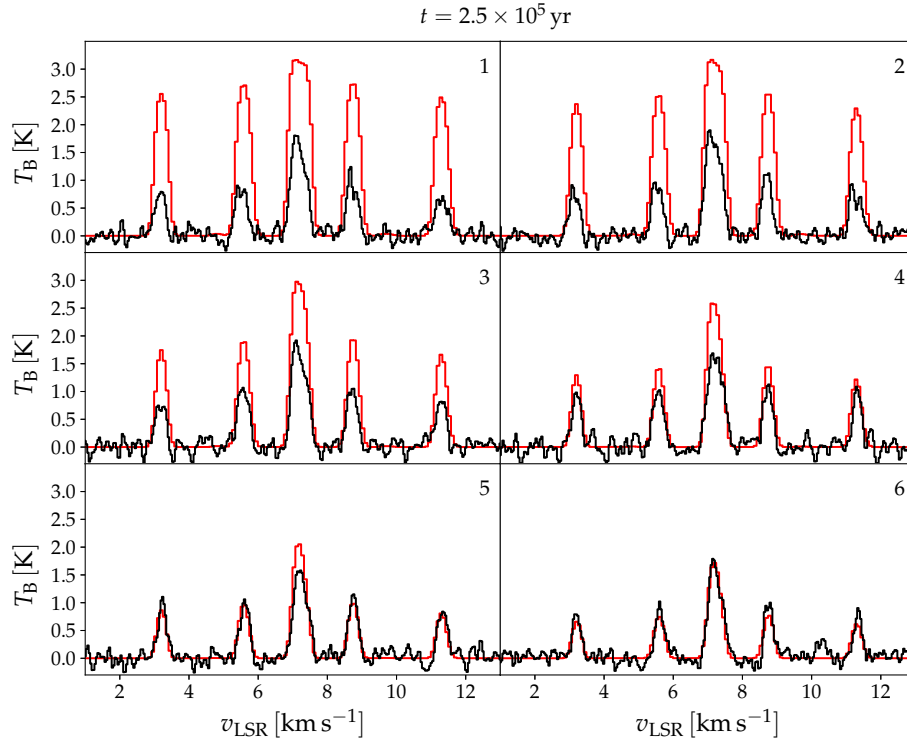


Figure C1. Same as Fig. 3, right panels, but assuming that the abundance of pNH_2D is constant within the central 2000 au, instead of dropping as predicted by our chemical models (see Fig. 5). The modeled lines clearly overestimate the observed spectra in the central four positions, demonstrating that freeze-out is needed.

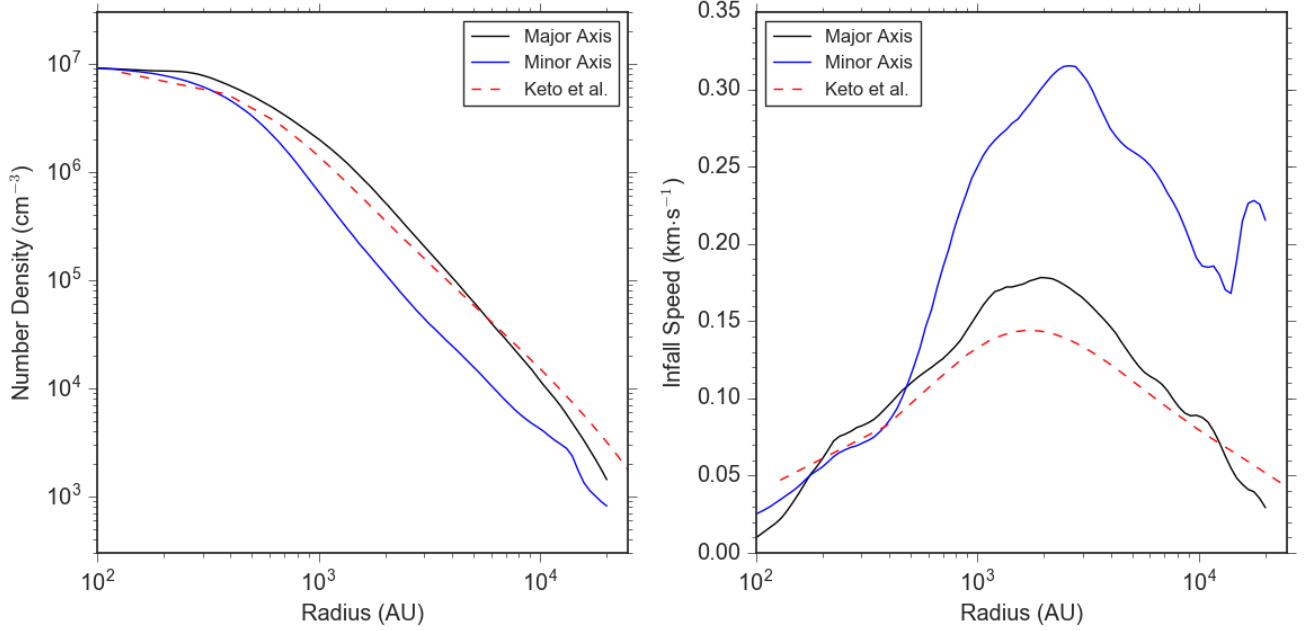


Figure D1. Comparison between the H_2 number density (*left panel*) and velocity (*right panel*) radial profiles of the BE sphere in quasi-static contraction from the HD simulations of Keto et al. (2015), after 6×10^5 yr of evolution of an initial BE sphere with central density $1 \times 10^4 \text{ cm}^{-3}$, radius $2 \times 10^5 \text{ au}$ and total mass $10 M_\odot$ (dashed red curves), and those along the major (black curves) and minor (blue curves) axis of the niMHD simulation of a magnetised contracting core, after 1.4×10^6 yr of evolution starting from a uniform $8.1 M_\odot$ cloud with volume density $2 \times 10^4 \text{ cm}^{-3}$, radius $5 \times 10^4 \text{ au}$ and initial mass-to-flux ratio of 1, which best fits the ALMA 1.3 mm dust continuum emission (see Caselli et al. 2019, for details). Note the overall similarities despite the very different simulations (HD vs. niMHD) and different initial conditions.

Despite the structural difference between the central 1000 au of the Keto et al. (2015) model (used as input of our chemical model) compared with the 3D niMHD simulated core (which better reproduces the ALMA 1.3 mm dust continuum emission), the agreement between model and observed spectra in Fig. 3 is excellent. This is partly due to the choice of the cut, taken along the (projected) major axis of the kernel, as this is the direction where the density profile of the 3D structure more closely resembles the BE sphere (see Fig. D1, left panel). The agreement with observed spectra along the minor axis becomes poorer (than that with observed spectra along the major axis) in the outer two positions, as seen in Fig. A2. To reproduce in detail the spectra along the minor axis, we will have to include the chemistry in the 3D niMHD simulation, and this will be done in a future paper. For now, it is interesting to point out that the BE sphere in quasi-static contraction of Keto et al. (2015) has a density profile just in between the minor and major axis of the 3D niMHD simulated core (Fig. D1, left panel). Also the radial velocity profile of the contracting BE sphere is very similar to that along the minor and major axis within a radius of 500 au (Fig. D1, right panel); above this radius, the velocity profile of the contracting BE sphere follows more closely that along the major axis of the 3D niMHD core and it is only a factor of 2 away from that along the minor axis. Considering the simplicity of the contracting BE sphere HD simulation compared with the niMHD simulation of the contracting magnetized core, it is striking to see how closely the two resemble each other.

E. MODEL PREDICTIONS ASSUMING A $5'' \times 5''$ BEAM

Crapsi et al. (2007) observed the same transition of deuterated ammonia using the IRAM Plateau de Bure interferometer with an angular resolution of $5.8'' \times 4.5''$. To demonstrate that this is not sufficient to detect the pNH_2D depletion zone, we perform simulated observations of the same modelled core described in Sect. 4, assuming a spherical half power beam width of $5''$. Fig. E1 presents six simulated spectra extracted from six positions similar to those shown in Fig. 3, but now separated by $5''$ (instead $2.5''$). A $5''$ beam does not have a substantial effect on the simulated lines toward the center of the core (when compared with the red spectra in Fig. 3). The most noticeable difference is that the satellite components become slightly less bright (and less optically thick) compared to the simulations with a $2.5''$ beam; the line brightness now peaks toward the central two positions (number 1 and 2 in Fig. E1). An integrated intensity map would then show a peak at the center and a monotonic drop toward the outer parts, consistent with the Crapsi et al. (2007) observations. Therefore, with $5''$ resolution, as for the integrated intensity map, the central flattening in the column density profile could not be detected, because it is only present within the central resolution elements of Crapsi et al. (2007) observations.

We would also like to point out another important difference between our ALMA observations and those carried out by Crapsi et al. (2007): our new ALMA image of $\text{pNH}_2\text{D}(1_{11} - 1_{01})$ is the result of a 3-point mosaic (see Fig. A1), while the IRAM-PdBI observations consist of a single-point image without including a primary beam correction. This may cause an artificially steeper flux drop away from the central region, mimicking a centrally concentrated pNH_2D flux profile consistent with the distribution of the dust continuum emission. Finally, this leads to the misleading conclusion that central freeze-out is not needed for deuterated ammonia, pointing out the importance of high-angular resolution and high-sensitivity mosaic observations for a correct interpretation of pre-stellar core observations.

F. $\text{NH}_3(1,1)$ SIMULATED SPECTRA

Given that previous work has claimed that the fractional abundance of NH_3 is actually increasing toward the L1544 center (e.g. Crapsi et al. 2007), we would like to test if our model predictions about the NH_3 fractional abundance across the core (see Fig. 5) is consistent with previous interferometric work. For this, we simulate VLA $\text{NH}_3(1,1)$ observations of our model cloud using the LOC radiative transfer code and assuming a telescope beam of $4''$ (close to the VLA synthesized beam in the observations of Crapsi et al. 2007, $4.34'' \times 3.45''$), at two different positions: the dust peak and the offset ($10''$, $-20''$), as in Figure 2 of Crapsi et al. (2007). The resultant spectra, focusing on the observed central three groups of hyperfine components, are displayed in Fig. F1 (solid histogram). Both spectra are in good agreement with the VLA observations, taking into account their 15% flux uncertainty. Also, we are not simulating interferometric-only observations, as reported in Crapsi et al. (2007), as we take into account all the flux from the various scales, including the extended envelope, filtered out by the VLA observations. This point will be made clear in Schmiedeke et al., in prep., where a higher-sensitivity map of $\text{NH}_3(1,1)$ line obtained by combining VLA with GBT data will be presented. Therefore, we conclude that our model, where NH_3 presents a sharp drop in abundance toward the pre-stellar core center, is consistent with VLA observations. We also considered a case where the NH_3 fractional abundance does not drop in the central 6000 au (unlike in our models, shown in Fig. 5) and instead it maintains a constant value of 2×10^{-8} . The resultant spectra are also in Fig. F1 (see dashed histograms); they are almost identical to the solid histograms, underlining the fact that $\text{NH}_3(1,1)$ data cannot be used to constrain the abundance profile of NH_3 within the central region of the core. This is due to the relatively low critical density of $\text{NH}_3(1,1)$ ($\simeq 4 \times 10^3 \text{ cm}^{-3}$; Maret et al. 2009), which then makes the 1,1 inversion transition mainly sensitive to the outer parts of the pre-stellar core, hiding abundance variations in the central few thousand au. In a sense, this result is similar to that found for the $\text{pNH}_2\text{D}(1_{11} - 1_{01})$ line, with the important difference that $\text{pNH}_2\text{D}(1_{11} - 1_{01})$ has a significantly higher critical density than $\text{NH}_3(1,1)$ and then it is more sensitive to variations in the NH_2D fractional abundance within the central regions (as demonstrated by Fig. C1).

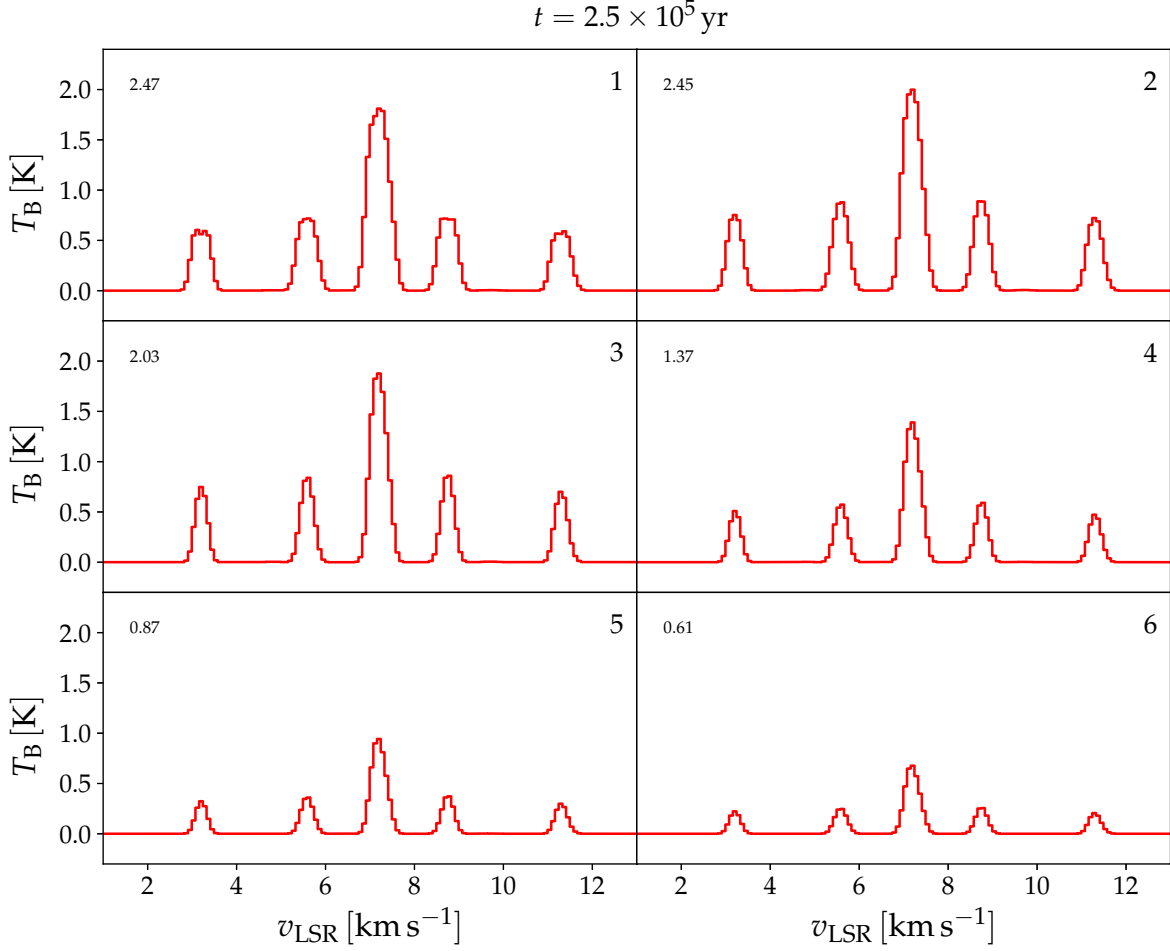


Figure E1. Simulated spectra of the same L1544 core model presented in Section 4 and Fig. 3 but now assuming a telescope beam of $5'' \times 5''$, similar to the one used by Crapsi et al. (2007). The number in the top-left of each panel is the integrated intensity in K km s^{-1} , while the number in the top right shows the corresponding position across the strip labelled in Fig. 3, left panel, but with a spacing of $5''$ instead of $2.5''$. Although the simulated spectra look similar to those in Fig. 3, the difference in integrated intensity between the spectrum extracted toward the core center (position 1) and the one $5''$ away (position 2, the integrated intensity peak) is very small. Thus, the flattening zone cannot be detected with the $\sim 5''$ beam of the Crapsi et al. (2007) observations.

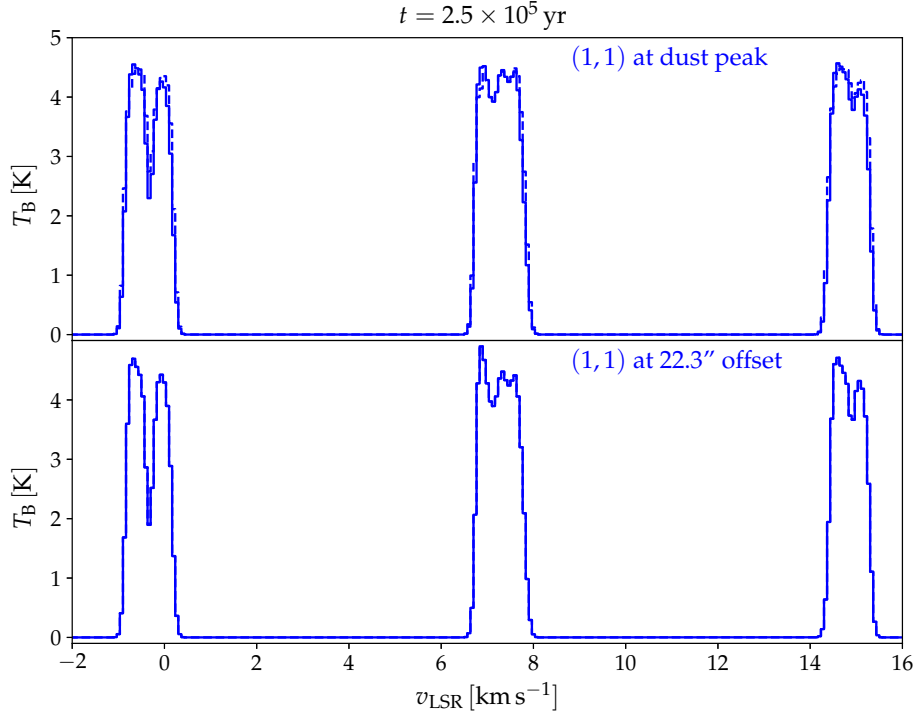


Figure F1. Simulated $\text{NH}_3(1,1)$ spectra toward the dust peak (top panel) and an offset $22.3''$ (bottom panel) away from the center, to be compared with the two spectra in Figure 2 of Crapsi et al. (2007). A beam of $4''$ (close to the angular resolution of the VLA observations of Crapsi et al. 2007) has been assumed. The solid blue histogram is the spectrum obtained with the NH_3 abundance profile predicted by our model and shown in Fig. 5, while the dashed blue histogram (only visible in the top panel, as in the bottom panel it coincides with the solid histogram) is the spectrum predicted assuming NH_3 constant abundance within the central 5000 au. Both spectra well reproduce the observed ones, within the observational uncertainties. It is then clear that $\text{NH}_3(1,1)$ interferometric spectra of L1544 cannot provide constraints on the fractional abundance distribution of ammonia, unlike ALMA observations of $\text{NH}_2\text{D}(1_{11} - 1_{01})$ (see Fig. C1).