# Temperature dependent ultracold neutron transmission in $D_2$ gas – a test of the Young-Koppel model

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The Young-Koppel model (YK) describes comprehensively the interaction of slow neutrons with diatomic gases such as  $H_2$  and  $D_2$ . This paper reports on the first experimental results of ultracold neutron (UCN) scattering over a wide temperature range vindicating the YK model for gaseous  $D_2$  and showing an important difference in the temperature dependence to a low-energy low-temperature approximation (LETA). LETA is confirmed, however, to be valid for monoatomic gases such as Ne. Calculated cross sections for other noble gases were also confirmed for ultracold neutrons. Finally, the total cross section of UCNs in  $H_2$  gas was measured and analyzed applying the Young-Koppel model, however, in a more limited temperature range, confirming the theoretical prediction.

# I. INTRODUCTION

Ultracold neutrons (UCN) are free neutrons with kinetic energies below about 300 neV that experience total reflection under all angles of incidence from surfaces of suitable materials [1]. The UCN source at the Paul Scherrer Institute [2–4] is used for various fundamental particle physics experiments, such as the search for an electric dipole moment of the neutron [5, 6] and a measurement of the free neutron lifetime [7].

The present work investigates the interaction between slow neutrons and individual deuterium molecules. This interaction, by its fundamental nature, has intrigued researchers from the early days of neutron physics [8–10] and remains highly topical until today in nuclear physics, for example, in the design of neutron sources [11]. It also has important practical implications for the nonstandard operating conditions of the PSI UCN source when considerable D<sub>2</sub> vapor pressure is present above the solid D<sub>2</sub> [12], as UCN-D<sub>2</sub> scattering can lead to significant UCN losses via up-scattering - i.e. neutrons acquire large enough energies to leave the UCN energy range.

In this work we tested the Young-Koppel model [13] that can be applied to describe the interaction between UCNs and  $H_2$  or  $D_2$  gases. It is the first time that the Young-Koppel model was tested over a wide temperature range from about 100 K to above 300 K.

# **II. THEORETICAL DESCRIPTION**

Young and Koppel (YK) calculated the neutron scattering cross sections from hydrogen and deuterium molecules by taking into consideration the spin correlation effects, as well as rotation and vibration of the molecules [13, 14]. In this approach, they considered vibrations as harmonic and decoupled from rotation. Also, there is no interaction between the molecules which would be the case in high-pressure gases or in liquids. The YK model can also be applied for liquids, however, for incident neutrons with energies less than roughly the solid state Debye temperature it is not accurate [13]. When the YK model was published, it was supported only by a limited amount of experimental data of the total cross section of neutrons in hydrogen gas as a function of incident energy [15].

Later, the hydrogen gas form factor was measured as a function of scattering angle with inelastic neutron scattering [16]. An excellent match was obtained with the YK model predictions, however, only in a very limited parameter space, at 18 K and one incident energy. The authors also reported a 20 % mismatch between the fitted and the observed temperatures, which they explained by a temperature gradient in the measurement cell.

A modified model for high pressures, assuming an effective temperature instead of the observed one, was tested in Ref. [17]. The measurements reported in [16] were completed in the same parameter space as reported in [18], in which the double-differential cross section of molecular hydrogen was measured at room temperature (RT) and could be fitted by using the YK model.

A significant time later, the total cross section of deuterium gas at 25 K was measured [19]. However, the YK model and a low energy approximation model could not be distinguished at this single low temperature, thus it did not provide a full corroboration of the YK model. New gas measurements were performed at room temperature [20, 21], where the authors scaled with  $\sqrt{T}$ the 25 K-value reported in Ref. [19] to an RT value. With simple reasoning, in the coordinate system of the molecules, the UCNs propagate with the mean velocity of the Maxwell-Boltzmann gas,  $w \propto \sqrt{T}$ , and the interaction rate  $\sigma w$  is invariant to changing the coordinate system (see also Eq. (4) in Ref. [19]). This assumption of the low-energy low-temperature approximation (LETA)

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FIG. 1. Ratio of the UCN upscattering cross sections calculated from the YK and LETA models revealing the discrepancy between the temperature dependences, and a local linear fit approximation with its parameters.

is over-simplified at higher  $D_2$  gas temperatures. The LETA model is similar to the monoatomic gas model [22], however, it uses an effective scattering cross section neglecting rotations and spin-correlation effects. The YK theory shows an important deviation of the up-scattering cross section from a pure  $\sqrt{T}$  dependence, which will be demonstrated experimentally in the present paper. Molecular vibrations are also included in the YK model, however, they are negligible in our experimental temperature range (vibration modes are populated only above 4000 K).

The total upscattering cross section as a function of UCN energies can be calculated from Eqs. (A4) and (A5) in [13] by integrating over the solid angle and final energies. We obtained a temperature dependence characteristic of the YK model. In Fig. 1 we plotted this dependence normalized with the LETA model, which only involves  $T^{1/2}$  dependence. At low temperatures, below the linearity range shown, the two model predictions co-incide, however, in the temperature range reaching above RT there is an important deviation. The YK model cross section is about 80% of that of LETA at RT.

#### III. EXPERIMENT

The experiment was designed to be sensitive to the difference between the monoatomic LETA and diatomic YK models, by measuring the transmission probability of UCNs passing through a gas as a function of pressure and temperature.

## A. Experimental setup

The experiment was conducted at beamport West-1 of the PSI UCN source, working with typically 8 s proton pulses on the target which start UCN measurements lasting for 300 s until the next pulse [2–4]. Beamport South was used to monitor the UCN source stability. We employed glass tubes coated with nickel-molybdenum (NiMo) and polished stainless steel tubes as UCN guides.



FIG. 2. The gas cell, chopper, and UCN detector (left to right, not to scale) in the standard TOF measurement. On the left side, the system was connected to the beamport West-1 of the UCN source.

The gas cell was followed by a UCN chopper [23], a 1 m UCN guide, and a CASCADE UCN detector [24] (see Fig. 2).

The gas cell was connected with the beamport to the left and a chopper connecting via a 1 m UCN guide to the UCN detector as in a standard time-of-flight setup. This chopper was only used for calibration measurements, otherwise left permanently open, thus ensuring stable conditions for the transmission measurements.

The gas cell was sealed on both sides with 0.1 mm thick aluminum (AlMg3) windows to separate the gas volume from the vacuum in the UCN guides. This window was supported by stainless-steel grids that covered approximately 10% of the tubes cross section. The gas cell was made of stainless steel as a compromise between temperature stability and UCN guiding properties. Cooling was done with a cold gas system where the cold  $N_2$  is guided first around the center of the cell and then in parallel to both ends. To achieve a good temperature equilibration with the windows, it was important that the cooling gas also cooled the stainless-steel UCN guides connected to both sides of the cell. Calculations and finite element simulations showed that the highest temperature gradient was expected starting from the center of the window to the point where the cold  $N_2$  gas is first reaching the cell. Therefore, we positioned several temperature sensors inside the gas volume which were introduced from the top via a pipe at a 45-degree angle to the horizontal. Additional temperature sensors were used on the outside of the cell on both aluminum windows. The T-sensors were calibrated beforehand with liquid N<sub>2</sub> and distilled ice water. In the final analysis, a time- and positionaveraged temperature value of the sensor readings was used. We also checked, that the impact of temperature gradients on the final result is negligible compared to our

main systematic uncertainty in the UCN dwell-time spectrum. Pressure sensors and the gas inlet were connected by a separate flange. We also made sure that there was no relevant effect of a temperature gradient on the pressure measurement.

## B. Analysis concept

The survival probability for a single UCN with a dwell time t and mean lifetime  $\tau$  inside the gas, when crossing the cell, is given by:

$$P = \exp(-t/\tau). \tag{1}$$

In an ideal system, the cell length, l, the velocity in axial direction,  $v_z$ , and the dwell time, t, are connected by the simple relation

$$t = l/v_z \tag{2}$$

The lifetime of the UCNs inside the gas cell is given by:

$$\tau = (\rho \ \sigma \ v_{\rm n})^{-1} = \left(\frac{p}{k_{\rm B}\sqrt{T}} \ \epsilon\right)^{-1},\tag{3}$$

where on the left-hand side,  $\rho$  is the number density of the gas molecules,  $v_n$  the neutron velocity and  $\sigma = \sigma(T, v_n)$ the microscopic up-scattering cross section, which is inversely proportional to  $v_n$  in both YK and LETA models. On the right-hand side, we used the ideal gas law (valid at our experimental conditions), where p is the pressure in the cell,  $k_B$  the Boltzmann constant, and introduced the parameter  $\epsilon$  defined as:

$$\epsilon = \frac{\sigma \, v_{\rm n}}{\sqrt{T}} \tag{4}$$

We introduced this scaled cross section parameter to make the analysis sensitive to temperature dependence other than  $\sqrt{T}$ . In case that LETA is valid,  $\epsilon$  would be independent of temperature. In case the YK model is valid,  $\epsilon$  would depend linearly on T (in the range of interest), see Fig. 1.

Putting everything together gives a  $v_z$ -dependent model for the analysis of the transmission probability

$$P_{\rm gas}^{v_z}(l, p, T; \frac{\epsilon}{v_z}) = \exp\left(-\frac{lp}{k_{\rm B}\sqrt{T}} \frac{\epsilon}{v_z}\right),\tag{5}$$

where the first three arguments describe the experimental conditions for the gas.

The quantity  $\epsilon$  is independent of temperature for monoatomic gases like Ne and is temperature dependent for diatomic gases like D<sub>2</sub>. As indicated in Fig. 1 for D<sub>2</sub>, the ratio of cross sections calculated with the YK and LETA models (thus also the ratio of the  $\epsilon$ -s defined via Eq. 4) can be approximated with a linear *T*-dependent fit function in the temperature range of interest (expecting discernible effects):

$$\epsilon(T) = \epsilon_0 \ (\epsilon_a - \epsilon_T \ T), \tag{6}$$

where  $\epsilon_0$  represents the LETA approximation, and  $\epsilon_a = 0.98 \pm 0.03$  and  $\epsilon_T = (6.26 \pm 0.37) \times 10^{-4} \text{ K}^{-1}$ . Besides the fit error, there is a ~ 1 % error from the numerical uncertainty of the MC integration and ~ 2 % from uncertainty in the input parameters like the equilibrium distance of the nuclei. In total, we estimate the slope error to be below 6 %.

Other than in the theory, in the analysis of the measurements, one cannot disentangle  $\epsilon_a$  from  $\epsilon_0$ . However, we were looking for the temperature dependence additional to  $\sqrt{T}$ . Thus, the signature we were searching for was covered by  $\epsilon_T$  and we thus set  $\epsilon_a$  to the theoretically expected value of 0.98 valid in our temperature range of interest.

In our analysis, we used calibrated (see next section) values for l, p (0 to 1 bar), T (range for all gases: 110 K to 365 K) and integrated Eq. 5 over the normalized neutron velocity distribution  $f(v_z)$  which was extracted from the combined time-of-flight (TOF) measurements and D<sub>2</sub> calibration (see Fig. 3 and details below) to obtain a measurable transmission probability

$$P_{\text{gas}}(l, p, T; \epsilon) = \int P_{\text{gas}}^{v_z}(l, p, T; \frac{\epsilon}{v_z}) f(v_z) dv_z, \quad (7)$$

where  $\epsilon$  also involves the parameter  $\epsilon_{\rm T}$ . As explained later, to also consider unavoidable UCN storage effects in the cell, integration over the velocity distribution was replaced in the analysis by the equivalent integration over dwell time (see Eq. 1).

We extensively investigated the measured data sets of two gases: neon (Ne) to serve as a test candidate for the LETA model and D<sub>2</sub> for the YK model. We checked whether  $\epsilon_T$  is compatible with 0 in the case of Ne, and whether for D<sub>2</sub> it is compatible with the value  $(6.26 \pm 0.37) \times 10^{-4} \,\mathrm{K^{-1}}$  from the theoretical Young-Koppel prediction.

As an additional check of systematics, mainly of the dwell time spectrum, we also measured He, Ar,  $CF_4$ , Xe and H<sub>2</sub> gases, but in a reduced p - T parameter space.

# **IV. SYSTEMATICS**

To extract the transmission probability function  $P_{\text{gas}}$ from the measurements, four major systematic effects had to be taken into account:

- time variation of the UCN rate at the beamport [25],
- faster neutrons during each beam pulse [2],
- temperature-dependent transmission probability of the empty cell, mainly of the windows,
- dwell time of UCNs in the gas cell.

A second CASCADE UCN detector (labeled here "B") on beamport South was used to monitor the UCN rate (measured with detector "A") at the West-1 beamport. Cutting the first 12s after the start of the proton beam pulse eliminated the contribution of neutrons faster than UCN [23]. In this way the ratio R between the neutron count rates of the two detectors for the empty cell, was determined to be constant at the < 1 % level.

The temperature-dependent transmission behavior of the empty cell was taken care of by calibration, as described in the second subsection below, applying a temperature-dependent correction factor for the total transmission probability.

Taking these three systematic effects into account  $P_{\text{gas}}(T)$  can be written as:

$$P_{\rm gas}(T) = \frac{1}{R \ P_{\rm window}(T)} \ \frac{N_A(p,T)}{N_B},\tag{8}$$

where  $N_A$  is the neutron count of the main detector and  $N_B$  is the neutron count of the monitoring detector for each proton beam pulse (both without the first 12 s).  $P_{\text{window}}(T)$  represents the temperature-dependent transmission probability of the empty cell and the UCN guides. The reference value R includes all other constant effects that influence the ratio between the number of counts in the detectors A and B.

# A. Reference value of the setup with empty cell

The reference R was measured periodically after and before thermal annealing of the D<sub>2</sub> crystal [25, 26] used for UCN production, for several UCN pulses with an empty cell. As the *T*-dependence of the UCN guides and cell without test-gas was reproducible, it could be done at any temperature  $T_{\rm ref}$  of the empty gas cell. However, most of the time it was measured at room temperature. The probability  $P_{\rm gas}(T)$  with an empty cell must be per definition 1, therefore, R can be written as:

$$R = \frac{N_{A,\text{ref}}(p=0,T_{\text{ref}})}{P_{\text{window}}(T_{\text{ref}})} \frac{1}{N_{B,\text{ref}}}$$
(9)

#### B. Temperature dependence for the empty cell

For the empty cell we found a thermal correction function by fitting the transmission data:

$$P_{\text{window}}(T) = \frac{C_T T + C_0}{C_0},\tag{10}$$

where  $C_T = (-1.73 \pm 0.03) \times 10^{-3} K^{-1}$  and  $C_0 = 5.470 \pm 0.006$ . Since it is impossible to disentangle the absolute transmission probability of the setup, the function was built in a way that it is 1 at 0 K. This is valid since the absolute scale will cancel out in  $P_{\rm gas}$  with the absolute scale of  $R(T_{\rm ref})$ . The relevant part for this investigation is, that the empty cell transmission changes

with the temperature so that the transmission probability of the empty cell at around 100 K is about 6 % higher than at RT. This is plausible considering up-scattering on phonons in the aluminum window causing a previously observed temperature dependence of the total UCN cross section on aluminum [27, 28].

#### C. Pressure-induced change of cell length

The effective length of the gas cell changes with increasing pressure in the cell, due to a small bulging of the aluminum foils on the grid. The size of this effect was estimated via a measurement of the pressure change in the gas cell while on the other sides of the windows the pressure changed from 1 bar to 0 during pumping. At a starting pressure of 972.0 mbar a small but well-measurable drop to 964.5 mbar was observed and could be reproduced several times. This was interpreted by ideal gas law as a linear change in the average cell length of 0.8 % from 0 to 1 bar with a linear dependence on the pressure.

Therefore, the effective mean length of the cell can be written as a function of pressure:

$$l = 127.6 \,\mathrm{mm} + p \times 1.060 \,\mathrm{mm/bar}$$
 (11)

This was in agreement with optical inspection and due to its small size a minor contribution in our analysis.

#### D. Calibration of the dwell-time spectrum

In our analysis the knowledge of the dwell-time distribution,  $N(t_{\text{dwell}})$ , i.e. the time UCNs spend inside the gas cell, is needed. In an ideal setup, where the axial velocity of the neutrons i.e. in the direction of the UCN guide axis,  $v_z$ , does not change over time,  $t_{\text{dwell}}$  can be derived from  $v_z$ :

$$t_{\rm dwell} = \frac{l_{\rm cell}}{v_z},\tag{12}$$

or  $v_z$  can be measured via time-of-flight (TOF) from the chopper to the detector:

$$v_z = \frac{l_{\rm TOF}}{t_{\rm TOF}}.$$
(13)

The UCN time-of-flight distribution was first measured with the chopper installed after the gas cell as in a standard TOF measurement (Fig. 3). However, this setup is blind to neutrons that are stored for a considerable time within the gas cell due to diffuse reflections, which causes an additional systematic.

To determine the size of this effect, we made use of the RT subset of all  $D_2$  transmission data measured with the chopper permanently open. First, from the TOF data taken with an empty cell, we performed an analytical



FIG. 3. The dwell time spectrum of UCNs is composed of two parts. The time axis was scaled using the ratio of the effective length of the gas cell and the flight path. First part: The time-of-flight spectrum of raw neutron counts measured with an empty cell (black dots 'TOF scaled'). The fit is denoted by ' $N_{\rm transmission}$ '. Second part: contribution of stored neutrons obtained from calibration to the RT data of D<sub>2</sub> denoted as ' $N_{\rm stored}$ '. In the analysis we used the combined spectrum ' $N_{\rm dwell}$ ' (see text).

parametrization of the main peak of the UCN spectrum. The profile can be described by a modified asymmetric Gaussian with a main peak around 0.025 s. This corresponds to a neutron velocity of 5.2 m/s (see TOF fit curve ' $N_{\text{transmission}}$ ' in Fig. 3). We checked for the impact of several factors like the precise shape of the main peak or small deviations at the rising and falling flanks, and found those to be negligible in our case.

As the second step, to describe the part of stored neutrons that are not visible in the TOF data we added an exponential tail:

$$N_{\text{stored}} = a_{storage} \exp\left(-t_{\text{dwell}}/\tau_{storage}\right) \qquad (14)$$

and on the left side a rising edge of the stored UCN curve which is described with a logistic growth function to emulate the filling of the system. The impact of the precise shape of the latter was found to be negligible. The final curve for the stored UCNs with calibrated parameters is shown as ' $N_{\text{stored}}$ ' in Fig. 3). The calibration of the exponential constants was determined by a global fit of Eq. 7 to all room-temperature D<sub>2</sub>-transmission data:

$$\tau_{storage} = 112.5 \pm 8.0 \text{ ms}, 
a_{storage} = 968 \pm 115.$$
(15)

Absolute amplitude values cancel out through later normalization, however, they are an indicator for statistical error estimation.

By this, we have a complete and calibrated description of the dwell-time spectrum. This was used in the analyses of all investigated gases at all temperatures and with both models (LETA and YK).

As an independent check of this exponential part, we rearranged the setup so that the chopper was placed before the cell. With this setup, we found a storage time constant of UCNs with around 92 ms, which is in reasonable agreement with the time constant, we found through the calibration procedure with  $D_2$  gas.

#### E. Systematic test with reference gases

To check our understanding of the dwell-time, we measured UCN transmission of gases as He, Ne, Ar,  $CF_4$ , Xe as references, and derived the total neutron cross section for a given gas at a given temperature.

Again, accounting for the correct dwell-time spectrum including stored UCNs we found good agreement with the literature and calculated the values shown in Table I. For reference, we considered 6.6 m/s UCN-velocity in order to compare to previous work [20] applying the same textbook equation (Eq. (11) in [20] from Ref. [29]) to convert from free atoms at rest to a Maxwell-Boltzmann gas. The values for the total cross section were obtained by fitting Eq. 7 to transmission data using the dwell-time spectrum from Fig. 3. The statistical uncertainty is extracted from the fit. The systematic uncertainty was obtained by maximum error propagation by fitting with minimum and maximum values derived from the uncertainties on  $\tau_{\rm storage}$  and  $a_{\rm storage}$  (Eq. 15).

By this, we find that for all investigated gases the systematic uncertainties dominate the statistical ones by at least one order of magnitude.

TABLE I. Measured total cross sections calculated for 6.6 m/s and at T=294.5 K including statistical and systematic uncertainties. Most 'literature' values were calculated from tabulated data [30]. Absorption data were all calculated from Ref. [30]. The CF<sub>4</sub> cross section is cited from ([20]). The 'literature' cross sections for D<sub>2</sub> and H<sub>2</sub> were calculated in this work by applying the YK model.

		literature			
species	σ	stat.	syst.	σ	
	$\mathbf{barn}$	$\pm$	±	barn	
He	185	1.13	17.0	164	
Ne	255	0.566	22.4	214	
Ar	287	1.74	24.9	264	
Xe	8220	34.1	389	8063	
$CF_4$	3450	1.87	141	3300	
$D_2$	2720	4.36	118	2582	
$H_2$	17400	49.7	969	18060	

# V. RESULTS AND DISCUSSION

The main transmission measurements were performed with gaseous  $D_2$  and Ne in the pressure range from 0 to 1 bar, and from 110 K to 365 K. The pressure, p, and the temperature, T, of the gas are measured directly with sensors in each measurement.

In Eq. 5 the exponent of the transmission probability of a neutron is given as  $\left(-\frac{lp}{k_{\rm B}\sqrt{T}} \frac{\epsilon}{v_z}\right)$ , where the cell length, l, is dependent on the absolute pressure, p, according to Eq. 11. Consequently, the transmission can be plotted as a function of  $\left(\frac{lp}{\sqrt{T}}\right)$ . By this choice, one expects a closeto-exponential dependence of the measured transmissions on  $\left(\frac{lp}{\sqrt{T}}\right)$ , which also includes the broadening due to the dwell time spectrum. These transmission profiles are seen in the top plots of Fig. 4 for D<sub>2</sub> gas and Fig. 5 for Ne.

The aforementioned spectral broadening would make calculating  $\epsilon(T)$  directly as a function of transmission (by inverting Eq. 7) to an arduous computational challenge. Thus we developed a feasible numerical fit alternative. In the data fit, we chose the transmission probability as the dependent variable, and the two independent variables were thus T and  $\left(\frac{lp}{\sqrt{T}}\right)$ , which replaced p. The unknown quantity given by the two-variable fit thus was  $\epsilon$ . Therefore, for the LETA model a single-parameter ( $\epsilon$ ) fit and for the YK model a two-parameter ( $\epsilon_0, \epsilon_T$ ) fit was performed. The latter additional parameter  $\epsilon_{\rm T}$  gave a small modulation to the exponential dependence, due to the additional T-dependent term, which served as the signature of the YK model. Our least-square fit minimized the sum of the transmission residuals from all measurements for a given gas at different T and  $\left(\frac{lp}{\sqrt{T}}\right)$  values. The residuals were defined here as the difference between the measured UCN transmission and the one calculated via Eq. 7, taking into account that  $\epsilon$  has the form as in Eq. 6.

In the top of Figs. 4 and 5, along with the measured transmissions, we also plotted the best fits with the two models for  $\epsilon$  plugged into Eq. 7 (corresponding to YK and LETA): one with  $\epsilon = \epsilon(T)$  (Eq. 6) and one with  $\epsilon$  independent of temperature. Since a full 3D plot of this two-variable fit problem would be much more difficult to interpret, we visualized the theoretical fit in the plots using only three example values for the independent variable T (100 K, 300 K and 400 K, not to be confounded with measured values).

In the case of the LETA model, the horizontal axis variable  $\left(\frac{lp}{\sqrt{T}}\right)$  is completely describing the  $\sqrt{T}$  dependence. In the case of the YK model, an additional temperature dependence occurs via the *T*-dependence of  $\epsilon$ . Thus for different *T*-s, we obtain different slopes for the exponent in the transmission versus  $\left(\frac{lp}{\sqrt{T}}\right)$  corresponding to each  $\epsilon(T)$ . This results in splitting the exponential transmission curve into separate curves corresponding to the number of plotted temperatures. Exactly this split-



FIG. 4. Top: UCN transmission data (black dots) in D<sub>2</sub> gas as a function of  $\left(\frac{lp}{\sqrt{T}}\right)$  (at all measured *T*-s), and the best fits of the two alternative models with example values for the independent variable *T* in case of YK (blue=100 K, yellow=300 K, red=400 K lines), and one line with LETA (black). Center: Point-by-point residuals of the YK and LETA model fit to the measured data, showing data at all measured temperatures in a color scale. Bottom: Histograms of the above residuals showing a much narrower and symmetrical distribution for the case of the YK model.

ting is the signal to distinguish between the LETA and YK models, as visualized in Fig. 4 for  $D_2$  and in Fig. 5 for Ne. Here we plotted three lines (i.e. three 3D-slices) for YK corresponding to three example values of the independent variable T enclosing the range of the measured temperatures.

In contrast to  $D_2$ , the Ne data fits with different temperatures shown in Fig. 5 converge with a slope in T compatible with zero, see also Table II, hence the fit involving Eq. 6 reproduces the LETA model.

The best fit of the Ne data results in  $\epsilon_T = (54 \pm 38 \pm 10) \times 10^{-6} \text{ K}^{-1}$ , which is about compatible with zero considering the uncertainties. For D<sub>2</sub> gas data the fit yielded  $\epsilon_T = (559 \pm 10 \pm 6) \times 10^{-6} \text{ K}^{-1}$ , with very high significance in terms of uncertainty.

The central plots display the residuals for both model fits. No distinction between the models is visible in the



FIG. 5. Top: UCN transmission data (black dots) in Ne gas as a function of  $\left(\frac{lp}{\sqrt{T}}\right)$  (at all measured *T*-s), and the best fits of the two alternative models with example values for the independent variable *T* in the case of YK (blue=100 K, yellow=300 K, red=400 K lines), and one line with LETA (black). The black and colored curves overlap (see insert), since no *T*-dependence of  $\epsilon$  was obtained from the Ne gas fit. Center: Residuals of the LETA and YK model fits to the measured data, here displaying all the measured temperatures in color scale. Bottom: Histograms of the above residuals.

Ne case, hence the parameter  $\epsilon_T$  does not contribute. In the D<sub>2</sub> gas case, this is clearly different: the LETA model shows systematic deviations that vanish when using the YK model. The same is true in the histogram of residuals in the bottom plots in Figs. 4 and 5.

The fit results are summarized in Table II. As in the results for other gases (Table I), the uncertainties of  $\epsilon_0$  are for both gases, D<sub>2</sub> and Ne, dominated by the systematic contribution from  $N_{\rm dwell}$ , whereas, for  $\epsilon_T$  the uncertainties are dominated by statistics, as the latter is a relative parameter. The neutron counts per measurement point for both detectors were above  $3 \times 10^5$ . However, considering only pure neutron count rate statistics would neglect other effects, such as minor changes in the neutron spectrum due to slow changes in the neutron production efficiency of the source.

The cross-check with Ne gas gives further confidence

that in our current investigation, the function  $N_{\text{dwell}}$  fitted to the extended tail, which represents stored UCNs, is appropriate, even if improvable for measuring absolute cross sections with higher accuracy.

TABLE II. Best model fit parameters for Ne and  $D_2$  data with an additional parameter for a linear temperature dependence of the up-scattering cross section according to the YK model as shown in Eq. 6 and Fig. 1.

	-		0				
species	$\epsilon_0$	stat.	syst.	$\epsilon_T$	stat.	syst.	effective $\epsilon$
	$\left(\frac{\operatorname{barn}m/s}{K^{1/2}}\right)$	$\pm$	$\pm$	$10^{-6}/K$	$\pm$	$\pm$	$\left(\frac{\operatorname{barn} m/s}{K^{1/2}}\right)$
Ne	100.6	1.0	10.0	54	38	10	97.0
$D_2$	1238.8	3.8	54.8	559	10	6	1006.3

We interpret the large  $\epsilon_T$  value in the case of D<sub>2</sub> as a clear demonstration of an additional *T*-dependent term in the YK-model compared to the LETA model for monoatomic gases. Comparing this result to the theoretical prediction  $\epsilon_T = (626 \pm 37) \times 10^{-6} \text{ K}^{-1}$ , and considering the measurement uncertainties and the unevenly distributed residuals (center plots in Figs. 4 and 5) hints towards yet unresolved systematic uncertainties in the measurements.

Returning to Table I, the absolute value of the  $D_2$  total cross section for UCNs is given under 'literature' as calculated by integrating Eqs. (A4) and (A5) in Ref. [13] and adding in a 2:1 ortho-para RT-equilibrium ratio. We also measured the total cross section of UCNs in H<sub>2</sub> gas. The data were analyzed by applying the Young-Koppel model, however, in a more limited temperature range (250-350 K) and thus with less sensitivity to the details of this model. The theoretical YK differential cross sections, Eqs. (29)-(30) in Ref. [13], for para- and orthohydrogen were integrated over final energy and solid angle, and added in a 1:3 para to ortho equilibrium ratio at RT. The measurement result for H<sub>2</sub> and the theoretical expectation are in very good agreement, see Table I, consistent within the systematic uncertainties.

# VI. SUMMARY AND CONCLUSION

We performed UCN transmission measurements in  $D_2$ and other gases at temperatures between 110 K and 365 K. A comparison between the molecular deuterium and the monoatomic neon gas showed a clearly different temperature dependence of the total cross section. We analyzed our data using the Young-Koppel model and a Low-Energy Low-Temperature Approximation. We interpret our best fit to the  $D_2$  gas data as a strong confirmation for an additional temperature-dependent term due to spin correlations and rotational molecular states as described by the Young-Koppel model. This is an important confirmation of the YK theory and has implications for the understanding of deuterium-based ultracold neutron sources.

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