Spectroscopy of atomic hydrogen

How is the Rydberg constant determined?

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Abstract. This article presents a review of the most recent theoretical and experimental results in hydrogen. We particularly emphasize the methods used to deduce the Rydberg constant $R_\infty$ and we consider the prospects for future improvements in the precision of $R_\infty$.

1 Introduction

The hydrogen atom has a central position in the history of 20th-century physics. As it is the simplest of atoms, it has played a key role in testing fundamental theories, and hydrogen spectroscopy is associated with successive advances in the understanding of atomic structure. The advent of tunable lasers and nonlinear techniques of Doppler free spectroscopy in the seventies led to major advances in resolution and measurement precision which are described in references [1,2]. Since, in the nineties, hydrogen spectroscopy has inspired a revolution in the art of measuring the frequency of light thanks to optical frequency combs [3]. Consequently, several optical frequencies of hydrogen are now known with a fractional accuracy better than $10^{-11}$, the most precise being the 1S-2S two photon transition which has been measured with a relative uncertainty of $1.4 \times 10^{-14}$ [4]. Thanks to these advances, the accuracy of the Rydberg constant $R_\infty$ has been improved by several orders of magnitude in three decades. This is illustrated in the figure 1 which clearly shows the improvements due to laser spectroscopy and optical frequency measurements. Nevertheless, during the last decade, there has been little progress with, for example, no important improvement between the $R_\infty$ values given by the last two adjustments of the fundamental constants in 2002 and 2006 [5,6]. In this review, we describe the analysis of the theoretical and experimental data used to deduce $R_\infty$ and we discuss the prospects for the future improvements in the accuracy of $R_\infty$.

2 Energy levels of atomic hydrogen: theoretical calculations

2.1 Theoretical background

Figure 2 shows the energy levels of hydrogen for different steps of the theory. The solution of the Schrödinger equation gives the same energy levels as the simple Bohr model. The energy $E_n$ depends only on the principal quantum number $n$:

$$E_n = -\frac{\hbar c R_\infty}{n^2}$$

(1)

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Fig. 1. Relative precision of the Rydberg constant from 1920 to the present.

The Rydberg constant $R_{\infty}$ is known as a function of the fine structure constant $\alpha$, of the velocity of light $c$, of the Planck constant $h$ and of the electron mass $m_e$:

$$R_{\infty} = \frac{\alpha^2 m_e c}{2h}$$  \hspace{1cm} (2)

This equation which links several fundamental constants is very useful for the adjustment of the fundamental constants. In particular, as $R_{\infty}$ is very well known, it is a link between the fine structure constant $\alpha$ and the $h/m_e$ ratio and it is a way to deduce $\alpha$ from the $h/m$ measurements \[7,8,9,10\].

The next step takes into account the relativistic corrections which are given by the Dirac equation. This equation lifts the degeneracy in $j$ ($j$ is the total angular momentum; $j = l \pm 1/2$) and explains the fine structure, but, for instance, the levels $2S_{1/2}$ and $2P_{1/2}$ are degenerate. This degeneracy disappears with the corrections due to quantum electrodynamics (QED), which are responsible for the Lamb shift between the $2S_{1/2}$ and $2P_{1/2}$ levels, first observed by Lamb and Retherford in 1947 \[11\]. In this paper, we wish to give only an idea of the theoretical calculations in hydrogen and more details are found in textbooks and recent review papers \[12,13,14\].

The hydrogen level energy can be conventionally expressed by the sum of three terms:

$$E_{n,l,j} = E_{n,l,j}^{\text{Dirac}} + E_{n,l}^{\text{Recoil}} + L_{n,l,j}$$  \hspace{1cm} (3)

where $E_{n,l,j}^{\text{Dirac}}$ is the energy given by the Dirac equation for a particle with a reduced mass $m_r = m_e (1 + m_e/m_p)^{-1}$ and $E_{n,l}^{\text{Recoil}}$ the first relativistic correction due to the recoil of the proton which depends only on the principal quantum number $n$. The last term $L_{n,l,j}$ is the Lamb shift which takes into account all the other corrections: the QED corrections, the other relativistic corrections due to the recoil of the proton and the effect of the proton charge distribution. The first two terms of equation (3) are exactly known as a function of the Rydberg constant $R_{\infty}$, of the fine structure constant $\alpha$ and of the electron to proton mass ratio $m_e/m_p$. In contrast, the calculation of the Lamb shift $L_{n,l,j}$ is a very difficult. It is obtained in a series of terms in powers of $\alpha$, $Z \alpha$ ($Z$ is the charge of the nucleus and $Z \alpha$ characterizes the interaction between the proton and the electron), $m_e/m_p$ and the root-mean-square charge radius of the
Fig. 2. Energy levels of atomic hydrogen for successive steps of the theory. The Dirac equation describes the fine structure splitting between the \( j = 1/2 \) and \( j = 3/2 \) levels and QED theory the splitting between the \( 2S_{1/2} \) and \( 2P_{1/2} \) levels.

proton \( r_p \). A very clear review of these different terms is given in the reference [5]. In this paper, we only present the most recent results.

2.2 Recoil corrections

The first recoil correction is:

\[
E_{\text{Recoil}}^n = -\frac{m_e^2c^2}{m_e + m_p} \left\{ \frac{(Z\alpha)^4}{8n^3} \right\}^n
\]

In terms of frequency, this first correction is about 24 MHz for the 1S level, i.e. \( 10^{-8} \) times the ionization energy of hydrogen. The following terms of the recoil correction vary as \( m_e^2c^2(m_e/m_p)(Z\alpha)^4 \) and \( m_e^2c^2(m_e/m_p)(Z\alpha)^5 \) and they have an exact expression. The next term, which varies as \( m_e^2c^2(m_e/m_p)(Z\alpha)6 \), has been calculated in the period 1988-1998 by several authors [15,17,18,19,20]. Table 1 summarizes the successive results which have been obtained and shows the difficulty of this kind of calculations. Finally, the value of this correction is now known with an uncertainty of 10 Hz for the 1S level.

2.3 Quantum electrodynamics corrections

The corrections due to QED are the main contribution to the Lamb shift. The self energy (SE) corresponds to the emission and reabsorption of virtual photons by the electron, and the vacuum polarization (VP) to the creation of virtual electron-positron pairs. A simple explanation of the self energy is given by the Welton model [21]. Because of the residual energy of the empty modes of the electromagnetic field (the energy \( \hbar\omega/2 \) of the harmonic oscillators), the electron is submitted to the fluctuations of the vacuum field which induce fluctuations in its position.
virtual photons. It varies as few Hz \cite{28,29}.

\[ Z\alpha \]

has been obtained numerically for the 1S level since the seventies \cite{22,23,24,25,26,27}. More recently, the one-loop correction to all orders in \( G \)

There are similar equations for the vacuum polarization correction. All the terms in equation 4 will be inserted by the editor

This effect modifies the Coulomb potential seen by the electron and is particularly important for the S levels: it reduces the binding energy, i.e. it increases slightly the energy of the S states \((l = 0)\), because, for the S states, the electron has a large probability

\[ \Psi(0) \]

\[ 2S_{1/2} \] and \( 2P_{1/2} \) levels (see Figure 2).

For the self energy, the lowest-order radiative correction (called ”one-loop” or ”one-photon” correction) is given by:

\[ F_{\text{SE}}^{(2)} = \frac{\alpha}{n^3} F(Z\alpha) m_e c^2 \] (5)

where \( F(Z\alpha) \) is a sum of terms in powers of \( Z\alpha \) and \( \ln(Z\alpha) \):

\[ F(Z\alpha) = A_{41} \ln(Z\alpha)^{-2} + A_{40} + A_{50}(Z\alpha) + A_{62}(Z\alpha)^2 \ln^2(Z\alpha)^{-2} + A_{61}(Z\alpha)^2 \ln(Z\alpha)^{-2} + G_{\text{SE}}(Z\alpha)(Z\alpha)^2 \] (6)

There are similar equations for the vacuum polarization correction. All the terms in equation 4 have an exact expression, except the last one \( G_{\text{SE}}(Z\alpha) \) which has been calculated intensively since the seventies \cite{22,23,24,25,26,27}. More recently, the one-loop correction to all orders in \( Z\alpha \) has been obtained numerically for the 1S, 2S, 3S and 4S levels with an uncertainty of a few Hz \cite{28,29}.

The following radiative correction takes into account the emission and reabsorption of two virtual photons. It varies as \( \alpha^2 \) and is given by:

\[ E^{(4)} = \left( \frac{\alpha}{\pi} \right)^2 \frac{(Z\alpha)^4}{n^3} F^{(4)}(Z\alpha)m_e c^2 \] (7)

where:

\[ F^{(4)}(Z\alpha) = B_{40} + B_{50}(Z\alpha) + B_{63}(Z\alpha)^2 \ln^3(Z\alpha)^{-2} + B_{62}(Z\alpha)^2 \ln^2(Z\alpha)^{-2} + B_{61}(Z\alpha)^2 \ln(Z\alpha)^{-2} + B_{60}(Z\alpha)^2 + \ldots \] (8)

The terms of equation 8 are calculated in references \cite{30,31,32,33,34}. However there is a controversy on the last calculated term \( B_{60} \). For the 1S level, Pachucki and Jentschura obtain \( B_{60} = -61.6(9.2) \) \cite{35} while Yerokin et al. give \( B_{60} = -127(38) \) \cite{36}. This difference corresponds to 6.6 kHz for the Lamb shift of the 1S level.

2.4 Other corrections

There are many of other terms in the calculation of the Lamb shift which are detailed in reference \cite{5}: the three-loop radiative corrections, the effect of the creation of virtual pairs \( \mu^+ \mu^- \) and \( \tau^+ \tau^- \), the radiative recoil corrections, the self energy and the polarization of the nucleus, and the effect of the non-zero size of the nucleus. This last term is important and is given for the proton by:

\[ E_{\text{NS}} = \frac{2}{3} \left( \frac{m_\tau}{m_e} \right)^3 \frac{(Z\alpha)^2}{n^3} m_e c^2 \left( \frac{2\pi Z\alpha r_p}{\lambda c} \right)^2 \] (9)
Table 2. Summary of the calculation of the 1s Lamb shift. The main uncertainties are due to the proton size correction (about 50 kHz) and to the two-loop corrections (about 3.3 kHz). The uncertainties in the one-loop correction (SE and VP) are essentially due to the uncertainty in the fine structure constant $\alpha$.

| Term of the Lamb shift                        | Value for the 1S level | Uncertainties |
|-----------------------------------------------|------------------------|---------------|
| Self-energy (one-loop)                        | 8 383 339.466 kHz      | 0.083 kHz     |
| Vacuum polarization (one-loop)                | −214 816.607 kHz       | 0.005 kHz     |
| Recoil corrections                            | 2 401.782 kHz          | 0.010 kHz     |
| Proton size                                   | 1 253 000 kHz          | 50 kHz        |
| Two-loop corrections                          | 731 000 kHz            | 3.300 kHz     |
| Radiative recoil corrections                  | −12.321 kHz            | 0.740 kHz     |
| Vacuum polarization (muon)                    | −5.068 kHz             | < 0.001 kHz   |
| Vacuum polarization (hadron)                  | −3.401 kHz             | 0.076 kHz     |
| Proton self-energy                            | 4.618 kHz              | 0.160 kHz     |
| Three-loop corrections                        | 1.800 kHz              | 1.000 kHz     |
| Nuclear size corrections to SE and VP         | −0.149 kHz             | 0.011 kHz     |
| Proton polarization                           | −0.070 kHz             | 0.013 kHz     |

1S Lamb shift 8 172 894(51) kHz

where $\lambda_C$ is the Compton wavelength of the electron. The rms charge radius of the proton $r_p$ is obtained from elastic electron-proton scattering experiments. There is a long history of analysis of these experiments. The most recent work gives $r_p = 0.895(18)$ fm which corresponds to a shift of the 1S level of about 1.2 MHz [37].

To summarize, we give in table 2 the value of the different terms contributing to the Lamb shift of the 1S level. For this calculation, we have used the value of $r_p$ from reference [37] and the values of $R_{\infty}$, $\alpha$ and $m_e/m_p$ given in reference [5]. The theoretical uncertainty in the 1S Lamb shift is estimated to be 3.7 kHz, mainly due to the calculation of the two-loop corrections. At this level, the precision is not limited by the uncertainties in $R_{\infty}$, $\alpha$ and $m_e/m_p$. On the contrary, the uncertainty due to the charge distribution of the proton $r_p$ is about 50 kHz, making it the largest source of uncertainty in the calculation of the 1S Lamb shift.

3 Precise measurements in hydrogen

In this section, we present the hydrogen frequency measurements which are used for the adjustment of the fundamental constants and for the determination of the Rydberg constant.

3.1 Lamb shift of the $2S_{1/2}$ level

Since the historic measurement of Lamb and Retherford [11], a number of measurements of the $2S_{1/2} - 2P_{1/2}$ splitting have been reported [38,39,40,41,42]. The recent results are shown in the figure 3. The most precise direct determination of this splitting is the one by Lundeen and Pipkin (1057.845(9) MHz [39]). The value obtained by Hagley and Pipkin in 1994 is an indirect determination deduced from the $2S_{1/2} - 2P_{3/2}$ splitting [10]. By using the theoretical value of the $2P_{1/2} - 2P_{3/2}$ fine structure splitting [26], this determination is 1057.842(12) MHz (see figure 3). The result of Drake in 1998 (1057.852(15) MHz [41]) is also an indirect determination, obtained by measuring the anisotropy of the emitted light in an applied electric field. The experiment of Pal’chikov, Sokolov and Yakovlev was performed by using atomic interferometry to measure the ratio between the $2S_{1/2}$ Lamb shift and the lifetime of the $2P_{1/2}$ level. This result is very precise (1057.8514(19) MHz [42]), but there is a controversy over the theoretical value of $2P_{1/2}$ lifetime which is used [43,44]. Finally, if we take into account only the direct measurements of the $2S_{1/2} - 2P_{1/2}$ and $2S_{1/2} - 2P_{3/2}$ splittings, we obtain a mean value of 1057.8439(72) MHz for the $2S_{1/2}$ Lamb shift.
3.2 The 1S-2S transition

The 1S – 2S transition is studied by Doppler-free two-photon spectroscopy, first proposed by Vasilenko et al. [45]. The principle is to place an atom in a standing wave formed by two counter-propagating laser beams with the same frequency. If the atom absorbs one photon from each beam, the total momentum of the absorbed photons is zero, and, consequently, there is a cancellation of the first order Doppler effect and of the recoil effect. The idea to apply this two-photon spectroscopy to the 1S-2S transition in hydrogen was immediately proposed by several authors, Cagnac et al. [46], Baklanov and Chebotayev [47] and Hänisch et al. [48]. It was a very attractive proposal: because of the very small natural width (1.3 Hz) of the 2S level, the quality factor of the 1S – 2S transition is about $2 \times 10^{15}$. However it was experimentally very difficult, due to the UV wavelength of this transition (243 nm) and the low transition probability.

Since the first observation by Hänisch et al. in 1975 [49], the 1S – 2S transition has been studied by several other groups in Southampton [50], Oxford [51,52,53], Yale [54] and MIT [55]. Nevertheless, the most important work has been performed by the group of Hänisch, who has continuously studied this transition since these first observations. In a long series of experiments, Hänisch has improved the precision on the measurement of the 1S – 2S frequency. The first determination of this frequency used an interferometric method with a calibrated absorption line of $^{130}$Te$_2$ [50,51,52,53]. In these experiments, the accuracy was limited by that of the $^{130}$Te$_2$ reference (about $2.7 \times 10^{-10}$). In the nineties, this limitation was overcome thanks to optical frequency measurements. In a first experiment, Hänisch used a frequency chain which linked the 1S-2S frequency (about 2466 THz) to a transportable CH$_4$-stabilized He-Ne frequency standard at 88 THz [58,59]. Now, this complex frequency chain has been superseded by a femtosecond laser frequency comb, which links in one fell swoop the Cs clock at 9 GHz to the optical frequency. Thanks to this technique, pioneered by Hall and Hänisch, Hänisch’s group has recently succeeded in measuring the 1S-2S interval with respect to a transportable Cs atomic fountain clock from the SYRTE (formerly Laboratoire Primaire du Temps et des Fréquences) in Paris [60]. This last measurement reduces the uncertainty to 34 Hz (i.e. a relative uncertainty of about $1.4 \times 10^{-14}$). The value obtained for the 1S – 2S frequency is:

$$\nu_{1S-2S} = 2 466 061 413 187 074 (34) \text{ Hz (10)}$$

This result is one of the most precise optical frequency measurements.

3.3 Two-photon spectroscopy of the 2S – nS and 2S – nD transitions

In Paris we began to study the 2S – nS and 2S – nD transitions in 1983. These experiments are complementary to those of the 1S – 2S measurements, because the Lamb shift of the 2S level has been measured precisely (see section 3.1), and, consequently, it is easy to extract the Rydberg constant from the 2S – nS/D interval. The two-photon 2S – nS/D transitions are induced in a metastable 2S atomic beam of hydrogen or deuterium collinear with the counter-propagating
laser beams. The excitation wavelength is in the near infrared, for example 778 nm for the 2S−8S/D transitions. The details of this experiment are described in references [61,62,63].

In our first experiments, we used an interferometric method to compare the hydrogen wavelengths to an iodine stabilized He-Ne laser. With this method, we determined the frequencies of the 2S−nD transitions in hydrogen and deuterium for the levels n = 8, 10 and 12 [64,65]. The relative accuracy was limited to 1.7 × 10−12 by the standard laser. In 1993, we measured the optical frequencies of the 2S1/2−8S1/2, 2S1/2−8D3/2 and 2S1/2−8D5/2 transitions in hydrogen with a frequency chain using two standard lasers (iodine stabilized and methane stabilized helium-neon lasers) and obtained a precision in the range of 10−11 [66,67]. More recently, we remade these measurements in hydrogen and deuterium with an accuracy better than one part in 1011 [68] using a new frequency chain with a new standard laser, namely a diode laser at 778 nm stabilized on the 5S-5D two-photon transition of rubidium. The frequency of this standard was measured with a frequency chain at the (LPTF) [69]. Finally, we completed these results by the measurement of the 2S−12S/D transitions [70]. A complete report of these experiments is given in reference [71]. For example, we obtain for the frequency of the 2S1/2−8D5/2 transition in hydrogen:

\[ \nu_{2S_1/2-8D_5/2} = 770 649 561.581.1 (5.9) \text{ kHz} \]  

and we have a similar result in deuterium. The relative uncertainty is limited to 7.6 × 10−12 because of the natural width of the 8D level (572 kHz) and the inhomogenous light shift experienced by the atoms passing through the Gaussian profile of the laser beams.

### 3.4 Frequency comparison between hydrogen frequencies

The goal of this method is to avoid absolute frequency measurements. The idea, proposed by Hänsch, is to compare the 1S−2S frequency with transitions whose energies are approximatively one-fourth that of the 1S−2S transition (see equation 1), such as the 2S−4P transition [49,72,53,54], or the 2S−4S/D two-photon transitions [73,74,75]. In our group, we have also performed an experiment based on the same idea, but this time by comparing the 1S−3S and the 2S−6S/D frequencies [76]. The three last experiments [54,75,76] provided a determination of the 1S Lamb shift with an uncertainty of about 50 kHz and these results are always used in the adjustment of the fundamental constants [50].

### 4 Determination of the Rydberg constant

The Rydberg constant is deduced from the data described in section 3 through a least squares adjustment. It is possible to make this adjustment with only the hydrogen data, the values of the fine structure constant α and the electron-to-proton mass ratio me/mp being given a priori [71], or to perform a global adjustment with the data concerning all the fundamental constants. Since 1998, the CODATA (Committee on Data for Science and Technology) has used the latter method to determine the value of \(R_\infty\). The value obtained in the 2006 CODATA adjustment is [6]:

\[ R_\infty = 10 973 731.568 527(73) \text{ m}^{-1} \]  

with a relative uncertainty of 6.6 × 10−12. The advantage of this method is to give the most accurate value, but the drawback is a mixing of all the experimental and theoretical results and it is difficult to see the most important input data. In the following sections, we describe several simple ways to determine the Rydberg constant.

#### 4.1 Determination of \( R_\infty \) from the 1S−2S interval

In equation 3, the Dirac energy \(E_n^{\text{Dirac}}\) and the recoil energy \(E_n^{\text{Recoil}}\) are exactly known as a function of the Rydberg constant. Thus, it is possible to rewrite this equation in the form:

\[ E_{n,l,j} = a_{n,j} \hbar c R_\infty + L_{n,l,j} \]
where \( a_{n,j} \) is a numerical coefficient which is an exactly known function of \( \alpha \) and \( m_e/m_p \), and whose value is approximatively given by the Bohr model \( a_{n,j} \approx -1/n^2 \) (see equation [1]). Then, the 1S – 2S frequency is:

\[
\nu_{1S-2S} = (a_{2,1/2} - a_{1,1/2})cR_\infty + (L_{2S_{1/2}} - L_{1S_{1/2}})/h \approx (3/4)cR_\infty + (L_{2S_{1/2}} - L_{1S_{1/2}})/h \]  

(14)

where \( L_{1S_{1/2}} \) and \( L_{2S_{1/2}} \) are the Lamb shifts of the 1S and 2S levels. In this equation, the frequency \( \nu_{1S-2S} \) is known with an uncertainty of 34 Hz, but the Lamb shift difference \( L_{2S_{1/2}} - L_{1S_{1/2}} \) is calculated with a precision of only about 44 kHz, because of the uncertainty in the proton radius \( r_p \). Consequently, in spite of the very high accuracy of the 1S – 2S measurement \((1.4 \times 10^{-14})\), the Rydberg constant can be deduced from equation [14] with a relative uncertainty of only \((1.8 \times 10^{-11})\).

4.2 Determination of \( R_\infty \) from the 2S – 8D interval

To deduce \( R_\infty \) from the 2S – 8D measurement, we can follow the same method. We have:

\[
\nu_{2S_{1/2}-8D_{5/2}} = (a_{8,5/2} - a_{2,1/2})cR_\infty + (L_{8D_{5/2}} - L_{2S_{1/2}})/h \approx \left( \frac{1}{4} - \frac{1}{64} \right) cR_\infty + (L_{8D_{5/2}} - L_{2S_{1/2}})/h \]  

(15)

In this expression, the uncertainty on the theoretical values of the Lamb shifts are respectively 2.5 Hz and 6.4 kHz for the 8D_{5/2} and 2S_{1/2} levels (because of the \( 1/n^3 \) scaling law of the Lamb shift, the uncertainty on the 2S_{1/2} Lamb shift is one eighth that of the 1S_{1/2}). Taking into account the experimental uncertainty \((5.9 \text{ kHz})\), see equation [14], we can finally extract \( R_\infty \) from equation [15] with a relative uncertainty of about \((1.1 \times 10^{-11})\).

Another way to obtain \( R_\infty \) is to combine the 2S_{1/2} – 8D_{5/2} frequency with the measurement of the 2S_{1/2} Lamb shift (see section 3.1). In this way, it is possible to eliminate the 2S_{1/2} level. Indeed, we have:

\[
\nu_{2P_{1/2}-2S_{1/2}} + \nu_{2S_{1/2}-8D_{5/2}} = (a_{8,5/2} - a_{2,1/2})cR_\infty + (L_{8D_{5/2}} - L_{2P_{1/2}})/h \]  

(16)

In this case, the theoretical uncertainties on the 2P_{1/2} and 8D_{5/2} Lamb shifts are very small \((80 \text{ and } 2.5 \text{ Hz})\) \(26\). With this method, the accuracy of \( R_\infty \) is limited by the uncertainties in the measurements of the frequencies \( \nu_{2P_{1/2}-2S_{1/2}} \) and \( \nu_{2S_{1/2}-8D_{5/2}} \). The relative accuracy of \( R_\infty \) is finally \(1.2 \times 10^{-11}\). This method is slightly less precise, but it does not use the experimental value for the rms charge radius of the proton.

4.3 Comparison of the 1S – 2S and 2S – 8D intervals

In the two preceding sections, we have seen that the accuracy of the Rydberg constant is limited by the theoretical or experimental uncertainties in the Lamb shifts. In fact, it is possible to avoid this difficulty by using the \( 1/n^3 \) scaling law for the Lamb shift. Numerous terms of the Lamb shift vary with the principal quantum number exactly as \( 1/n^3 \) (for instance the effect of the charge distribution of the nucleus), and the deviation from this scaling law has been calculated precisely by Karshenboim \(27\). He obtains for the 1S and 2S Lamb shifts:

\[
\Delta_2 = (L_{1S_{1/2}} - 8L_{2S_{1/2}})/h = -187.232 \text{ (5 MHz)} \]  

(17)

A more recent calculation of \( \Delta_2 \) is given in reference \(28\). If we combine this equation with equations [14] and [15], we obtain a system of 3 equations where the unknowns are \( R_\infty \), \( L_{1S_{1/2}} \) and \( L_{2S_{1/2}} \). We can then form a linear combination to eliminate the 1S and 2S Lamb shifts:

\[
7\nu_{2S_{1/2}-8D_{5/2}} - \nu_{1S_{1/2}-2S_{1/2}} \approx \left( \frac{57}{64} \right) cR_\infty + 7L_{8D_{5/2}}/h + \Delta_2 \]  

(18)
In this expression, the main uncertainty is due to the measurement of the $2S_{1/2} - 8D_{5/2}$ frequency ($7 \times 5.9$ kHz) and the Rydberg constant is deduced with a relative uncertainty of $1.4 \times 10^{-11}$. The advantage of this method is to deduce $R_{\infty}$ without the measurements of the $2S_{1/2}$ Lamb shift and of the proton radius $r_p$. Moreover, this technique is applicable to both hydrogen and deuterium (there is no precise measurement of the $2S_{1/2}$ Lamb shift in deuterium). For instance, from the measurements of the $1S_{1/2} - 2S_{1/2}$, $2S_{1/2} - 8D_{5/2}$ and $2S_{1/2} - 12D_{5/2}$ frequencies in hydrogen and deuterium, one obtains a value of $R_{\infty}$ with a relative uncertainty better than $10^{-11}$.

\[ R_{\infty} = 10 \, 973 \, 731.568 \, 54(10) \, \text{m}^{-1} \]  

Moreover, we can deduce the Lamb shift of the $1S_{1/2}$ and $2S_{1/2}$ levels in hydrogen:

\[ L_{1S_{1/2}}/h = 8 \, 172.837(26) \, \text{MHz} \]  

\[ (L_{2S_{1/2}} - L_{2P_{1/2}})/h = 1 \, 057.8447(34) \, \text{MHz} \]  

In this last result, we have used the theoretical value of the $2P_{1/2}$ Lamb shift ($-12.835\,99 \, (8)$ MHz [26]). This value of the $2S_{1/2}$ Lamb shift deduced from the optical frequency measurements is more precise than the direct determinations made by microwave spectroscopy.

Finally, if we take into account the theoretical calculations of the Lamb shift (see section 2), we can deduce a value of the rms charge radius of the proton ($r_p = 0.8765(80) \, \text{fm}$) which is more precise than the one deduced from the electron-proton scattering experiments. For this calculation, we have used $B_{60} = -94.3$, which is the mean value of the results of the references [35,36] (see section 2.3).

5 Conclusion

In conclusion, we have presented several methods to determine the Rydberg constant. There are several limitations, mainly the uncertainty in the rms charge radius of the proton and the uncertainties in the measurements of the $2S - nS/D$ frequencies in hydrogen. To reduce the first limitation, a precise measurement of $r_p$ is ongoing at the Paul Scherrer Institute by spectroscopy of muonic hydrogen. The principle is to measure the $2S - 2P$ energy difference in $\mu^-p$ by infrared spectroscopy [79]. In muonic hydrogen, the muon is very close to the proton because its mass is about 207 times that of the electron. Consequently, the effect of the proton charge distribution is about 0.93 THz for the $2S$ level of muonic hydrogen whereas it is only 146 kHz in hydrogen. The details of this project are given in references [80,81]. The goal is to obtain a relative accuracy of $10^{-3}$. Then the effect of the proton size for the $1S$ level of hydrogen would be known with an uncertainty of 2.5 kHz. Using equation 14, it will be possible to deduce $R_{\infty}$ with a relative precision of about $2 \times 10^{-12}$.

An alternative way to improve the precision of the Rydberg constant is to measure another optical frequency in hydrogen very precisely. Several groups are working in this direction. At the National Physical Laboratory, Flowers and colleagues have built a new experiment to improve the measurements of the $2S - nS/D$ hydrogen frequencies by using a femtosecond frequency comb [82]. In Paris, we intend to measure the optical frequency of the $1S - 3S$ two-photon transition [83,84]. This transition is also being studied by the group of H"{a}nsch. The same group is also working towards an experiment on the He$^+$ ion. All these efforts are promising to improve the accuracy on the Rydberg constant in the near future.

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