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Collision Processes of Li³⁺ with Atomic Hydrogen: Cross Section Database

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Abstract

Using the available experimental and theoretical data, as well as established cross section scaling relationships, a cross section database for excitation, ionization and charge exchange in collisions of Li³⁺ ion with ground state and excited hydrogen atoms has been generated. The critically assessed cross sections are represented by analytic fit functions that have correct asymptotic behavior both at low and high collision energies.

The derived cross sections are also presented in graphical form.

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1. Introduction

Collision processes of Li³⁺ ions with atomic hydrogen are of interest in magnetic fusion energy research in the context of TESPEL based CXRS plasma diagnostic and transport studies [1,2] and proposed liquid lithium plasma facing divertor concepts [3,4].

Processes of interest include:

estimated accuracy.

$$\text{Li}^{3+} + \text{H}(n_0 \ge 1) \to \text{Li}^{3+} + \text{H}(n > n_0),$$
 (1)

$$\text{Li}^{3+} + \text{H}(n_0 \ge 1) \to \text{Li}^{3+} + \text{H}^{+} + \text{e},$$
 (2)

$$\text{Li}^{3+} + \text{H}(n_0 \ge 1) \to \text{Li}^{2+}(n'\ell') + \text{H}^+,$$
 (3)

where $n(n_0)$ and ℓ are the usual quantum numbers of electronic states. In this report we present a complete set of cross sections for the processes (1)-(3). As basis for this cross section set we have used the available experimental and theoretical data from the literature, critically assessed for their accuracy, and appropriate, well-established cross section scaling relationships. Most of the available data are pertinent to the initial $n_0\ell_0=1$ s state of hydrogen atom. The data sources for each of processes (1)-(3), and the method (procedure) of evaluation or derivation of the cross section for each particular reaction, will be discussed in detail in the sections that follow. For each cross section we shall provide also information about its

The evaluated (or derived) cross sections of individual reactions within processes (1)-(3) for the lower initial levels n_0 are presented by analytic fit functions that have correct physical behavior at both low and high collision energies. For the cross sections with high initial n_0 , appropriate scaling relations are used.

2. Excitation

2.1. Excitation from the ground state ($n_0 \ell_0 = 1s$)

Cross section measurements for Li³⁺ induced 1s $\rightarrow n$ (≥ 2) transitions do not exist in the literature. There are two sets of cross section calculations by the

adiabatic superpromotion model for $1s \rightarrow n=2$ [5] and by the atomic orbital close coupling (AOCC) model (with 24 AOs) for $1s \rightarrow n=2,3$ [6]. There are also cross section calculations for $1s \rightarrow n$ (≤ 4) induced by Aq+ fully stripped ions with q=2,4,6,8,14,and 26 performed by two-center atomic orbital close-coupling (2C-AOCC) and classical trajectory Monte Carlo (CTMC) method [7,8] in the intermediate to high energy range ($\sim 10 \text{keV/amu} - 800 \text{keV/amu}$), and "hidden-crossings" superpromotion method [9,10] (for q=2,4 only) at low collision energies ($\leq 25 \text{ keV/amu}$). (More detailed information on this and other work is given in [11].)

The results of these extensive calculations show that for a given $1s \rightarrow n$ transition the excitation cross section $\sigma_{exc}^{(q)}(n)$ scales with q, for $q \ge 2$, as [11]

$$\sigma_n^{(q)} - q \chi(q) \sigma_n^{(2)}(E/q), \quad E/q \ge 10 \text{ keV/amu}$$
 (4a)

$$\sigma_n^{(q)} \sim q \chi(q) \sigma_n^{(2)}(E), \qquad E/q \le 10 \text{ keV/amu}$$
 (4b)

where
$$\chi(q)=2^{0.5238(1-\sqrt{\frac{2}{q}})}$$
.

2.1.1. Transitions to n = 2,3,4 levels

The cross sections for $1s \rightarrow n=2,3,4$ transitions in H by Li³⁺ impact have been determined by using the data from [5] and [6] (for n=2,3) and scaling the corresponding recommended cross sections for He²⁺ and Be⁴⁺ impact [11] according to the relations (4) and, then taking the average of obtained scaled cross sections. The obtained excitation cross sections are fitted by the analytic expression (n=2,3,4)

$$\sigma_n^{(q)}(n) = a_1 \left[\frac{\exp(-a_2/E)\ell n(1+a_3E)}{E} + \frac{a_4 \exp(-a_5E)}{E^{a_6}} + \frac{a_7 \exp(-a_8/E)}{1+a_9E^{a_{10}}} \right] \left(\times 10^{-16} cm^2 \right)$$
(5)

where E is expressed in keV/amu units, and the values of fitting parameters a_i are given in Table 1. The rms deviation of the fits is smaller than 1% for all n. Comparison between the fitting line and the theoretical results of Ref.

[5,6] is also displayed in Figure 1 for 1s·n=2 excitation. We note that $\sigma_{exc}(n=2)$ given by Esq. (5) agrees well with the results of superpromotion model [5] for $E \le 25 \text{ keV/amu}$. The accuracy of the cross sections given by Eq. (5) and Table 1 should be close to the accuracy of corresponding cross sections for He^{2+ and} Be⁴⁺ impact, from which they have been derived. For $E \le 20 \text{ keV/amu}$, their uncertainty (defined as combined experimental and/or theoretical error) is 30-80%, for $20 \le E(\text{keV/amu}) \le 200$ it is 20-30%, and for $E \ge 200 \text{ keV/amu}$ it is expected to be 15-20%, or better.

2.1.2. Transitions to n = 5.6 levels

The cross sections for the transitions $1s \rightarrow n=5,6$ have been determined by using the corresponding He²⁺ impact data from [11] and the scaling relations (4). These cross sections have somewhat simpler structure and can be fitted by the expression (n=5,6)

$$\sigma_{exc}(n) = b_1 \left[\frac{\exp(-b_2/E)\ell n(1+b_3E)}{E} + \frac{b_4 \exp(-b_5E)}{E^{b_6} + b_7 E^{b_8}} \right] \quad (\times 10^{-16} cm^2)$$
 (6)

where E is in keV/amu units. The values of fitting parameters b_i in Eq. (6) are given in Table 2. The rms deviation of the fits is smaller than 1%. The expected uncertainty of the cross sections given by Eq. (6) and Table 2 is $40 \cdot 100\%$ for E $\leq 5 \text{keV/amu}$, $30 \cdot 40\%$ for $5 \leq E$ (keV/amu) ≤ 20 , $30 \cdot 40\%$ for $20 \leq E$ (keV/amu) ≤ 200 , and better than 20% for E $\geq 200 \text{ keV/amu}$.

2.1.3. Transitions to $n \geq 7$ levels

The cross sections for $1s \rightarrow n \geq 7$ excitation transitions can be obtained by using the well-known scaling relation (see, e.g., [11])

$$\sigma_{exc}(n \ge 7) = \left(\frac{6}{n}\right)^3 \sigma_{exc}(1s \to n = 6),\tag{7}$$

which should be accurate for $E \ge 20 \text{ keV/amu}$. Given the large uncertainty (>30%) of σ_{exc} (1s \rightarrow n =6) cross section for $E \le 20 \text{ keV/amu}$, the scaling

formula (7) can be safely used also for E ≤ 20keV/amu.

2.2. Excitation from $n_0 \ge 2$ levels

Specific cross section measurements or calculations for transitions from the $n_0 \geq 2$ levels of H by Li³⁺ impact do not exist in the literature. Calculations exist only for the transitions from $n_0\ell_0=2$ s to higher $n\ell$ states, performed by the superpromotion model in the low energy region [12]. The contribution from the $n_0\ell_0=2$ p state to the total excitation cross section from the $n_0=2$ level to any higher level is expected to be significant. Total excitation cross section from the levels $n_0=2$ and $n_0=3$, however, do exist for He²⁺[9] and Be⁴⁺[10] impact. In determining the Li³⁺ impact excitation cross sections from $n_0=2,3$ we shall use these data.

2.2.1. Excitation from $n_0 = 2$ level

a) Transitions to n = 3,4,5 levels

The cross sections for transitions $n_0 = 2 \rightarrow n = 3,4,5$ have been determined by scaling the corresponding cross section data for He²⁺ and Be⁴⁺ impact from [11] according to Eqs. (4), and taking their average. The obtained cross sections have been then fitted to the expression (n=3,4,5)

$$\sigma_{exc}(2 \to n) = c_1 \left[\frac{\exp(-c_2/E)\ell n(1 + c_3 E)}{E} + \frac{c_4 \exp(-c_5 E)}{E^{c_6}} \right] \quad (\times 10^{-16} cm^2)$$
 (8)

with collision energy E expressed in keV/amu. The values of fitting parameters c_i are given in Table 3, and the rms deviation of the fit is well below 1%. The uncertainty of the data is expected to be 40-80% for E \leq 5keV/amu, 30-40% for 5 \leq E(keV/amu) \leq 20, 20-30% for 20 \leq E(keV/amu) \leq 200 and 15-20% for E \geq 200 keV/amu.

b) Transitions to $n \ge 6$ levels

The cross sections for $n_0 = 2 \rightarrow n = 6.10$ transitions have been determined by scaling the corresponding cross sections for He²⁺ (or proton) impact from Ref.

[11] according to Eq. (4a). The obtained cross sections can be expressed in the form

$$\sigma_{exc}(2 \to n = 6 - 10) = c_{2-n}\sigma_{exc}(2 \to 5)$$
 (9)

where coefficients c_{2-n} have the values given in Table 4. For the transitions $n_0 = 2 \rightarrow n \ge 11$, one can use the n^{-3} -scaling,

$$\sigma_{exc}(2 \to n) = \left(\frac{10}{n}\right)^3 \sigma_{exc}(2 \to 10) , n > 10$$
 (10)

The uncertainty of the cross sections given by Eqs. (9) and (10) should be by 5-10% higher than that for $\sigma_{exc}(2 \rightarrow 5)$.

2.2.2. Transitions from $n_0 = 3$

a) Transitions to n = 4,5,6 levels

For the $n_0 = 3 \rightarrow n = 4,5,6$ transitions in H, induced by Li³⁺ impact, the cross sections have been determined as an average of the corresponding He²⁺ and Be⁴⁺ impact cross sections of Ref. [11], scaled according Eq. (4a). So derived cross sections were then fitted to an expression having the same form as Eq. (8). The values of fitting parameters c_i are given in Table 5. The uncertainty of derived cross sections is 40-80% for E \leq 5keV/amu, 30-40% for $5 \leq$ E(keV/amu) \leq 100, and 20-30% for E \geq 100keV/amu.

b) Transitions to $n \geq 7$ levels

The cross sections for $n_0 = 3 \rightarrow n = 7 \cdot 10$ transitions have been determined by scaling the corresponding He²⁺ impact cross sections of Ref. [11] according to Eq. (4a). The obtained cross sections can be expressed in the form

$$\sigma_{eyc}(3 \to n = 7 - 10) = c_{3-n}\sigma_{eyc}(3 \to 6)$$
 (11)

where the coefficients c_{3-n} for n=7,8,9,10 have the values given in Table 6. For the transitions to $n \ge 11$ levels, the cross sections can be evaluated by using the n^{-3} -scaling

$$\sigma_{exc}(3 \to n) = \left(\frac{10}{n}\right)^3 \sigma_{exc}(3 \to 10), \quad n > 10$$
(12)

The uncertainty of the cross sections given by Eqs. (11) and (12) is expected to be by 5-10% higher than that for $\sigma_{exc}(3 \rightarrow 6)$.

2.2.3. Transitions from $n_0 \ge 4$

In absence of any systematic cross section data for transitions from $n_0 \ge 4$ levels, the q-scaled proton-impact semi-empirical formula of Lodge et al. [13] can be used. In order to account for the small deviations of the q-scaling at large energies and for high ionic charge q, a correction factor $\xi(q)$ has been introduced in Ref. [11]

$$\xi(q) = 2^{0.322(1-\sqrt{2/q})} \tag{13}$$

which for q = 3 is $\xi(3) = 1.042$. The q-scaled Lodge et al. formula [11] for q=3 and the $n_0 \to n$ transition can be written as

$$\sigma_{exc}(n_0 \to n) = 2.769 \frac{n_0^4}{\varepsilon} [ADL + FGH] \qquad (\times 10^{-16} cm^2)$$
 (14)

where

$$\varepsilon = \frac{E(keV/amu)}{75}, \quad s = n - n_0, \quad D = \exp\left[-\frac{1}{n_0 n \varepsilon^2}\right],$$

$$A = \frac{8}{3s} \left(\frac{n}{sn_0}\right)^3 \left(0.184 - \frac{0.04}{s^{2/3}}\right) \left(1 - \frac{0.2s}{n_0 n}\right)^{1+2s}, \quad G = \frac{1}{2} \left(\frac{\varepsilon n_0^2}{n - 1/n}\right)^3,$$

$$L = \ln \left(\frac{1 + 0.53\varepsilon^{2} n_{0} (n - 2/n)}{1 + 0.4\varepsilon} \right), \quad F = \left(1 - \frac{0.3sD}{n_{0} n} \right)^{1 + 2s},$$

$$H = \left[c_{2} \left(z^{(-)}, y \right) - c_{2} \left(z^{(+)}, y \right) \right], \quad c_{2} \left(z, y \right) = \frac{z^{2} \ln \left(1 + 2z/3 \right)}{2y + 3z/2},$$

$$z^{(\pm)} = \frac{2}{\varepsilon n_{0}^{2} \left[\left(2 - \frac{n_{0}^{2}}{n^{2}} \right)^{\frac{1}{2}} \pm 1 \right]}, \quad y = \frac{1}{1 - D \ln \left(18s \right) / 4s}.$$

The accuracy of the cross sections given by Eq. (14) is expected to be by 5·10% worse than for the $\sigma_{exc}(2 \to 5)$ or $\sigma_{exc}(3 \to 6)$ cross sections.

3. Ionization

3.1. Ionization from the ground state ($n_0 \ell_0 = 1s$)

The ionization of H(1s) by Li³⁺ impact has been subject to extensive both experimental and theoretical studies. The experimental cross section measurements exist in the energy range 50 - 500 keV/amu [14,15] and have accuracy on the level 10%. Theoretical cross section calculations have been performed by using the adiabatic superpromotion model [5], atomic-orbital close-coupling (AOCC) method [6,16], molecular-orbital close-coupling (MOCC) method [17], classical trajectory Monte Carlo (CTMC) method [18] and continuum distorted wave (CDW) method [19]. These results are plotted in figure 2, as well as the fitting line using the formula described below. The theoretical calculations cover the energy range from $\sim 1 \text{ keV/amu}$ to ~ 650 keV/amu, and agree with each other within 10-20% in the overlapping regions of their validity. While the MOCC method is valid at energies below valid down to $\sim 1-5$ keV/amu. The validity of CTMC methods is restricted to the range ~30-200 keV/amu, while CDW is a typical high-energy method valid for E ≥ 150-200 keV/amu. For E > 200-300 keV/amu, the Bethe-Born approximation (BBA) should also be valid. The BBA gives the ionization cross section for H(1s) in closed form for any fully stripped ion [20] (q-scaled cross section). It should be noted that the low-energy ionization cross section for H(1s) by fully stripped ion impact allows also a q-scaled representation [21]. The Li³⁺+H(1s) ionization cross section recommended for use is based on the very extensive AOCC calculations [16], experimental data [14,15], and CDW and q-scaled BBA cross sections. The AOCC cross section for E \leq 20 keV/amu is consistent with the cross section derived from the adiabatic superpromotion data [5] and q-scaling formula [21]. So determined "recommended" ionization cross section for H(1s) by Li³⁺ impact was fitted to the analytic expression ($n_0 = 1$)

$$\sigma_{ion}(n_0) = a_1 \left[\frac{\exp(-a_2/E)\ell n(1+a_3E)}{E} + \frac{a_4 \exp(-a_5E)}{E^{a_6} + a_7 E^{a_8}} \right] \qquad (\times 10^{-16} cm^2)$$
 (15)

where the collision energy is in keV/amu units and the values of fitting parameters are given in Table 7. The rms deviation of the fit is about 5%, most of it generated in the range 7-30 keV/amu where the AOCC cross section data are about 10-20% below the fit. By choosing a more complicated fit expression for $\sigma_{ion}(ls)$ (such as given by Eq. (5), for instance), the rms could be reduced to a much lower level. However, for the sake of uniformity with fitting expressions for $\sigma_{ion}(n0)$, for $n_0=2,3$, (see next subsection), we have kept the fit of $\sigma_{ion}(ls)$ in the form of Eq. (15). The estimated uncertainty of $\sigma_{ion}(ls)$ is 15-20% for E \leq 50 keV/amu and 10-15% for E \geq 50 keV/amu.

3.2. Ionization from $n_0 \ge 2$ levels

3.2.1. Ionization from $n_0 = 2$ and $n_0 = 3$

There are no cross section data for Li³⁺ impact ionization of $H(n_0)$ for the n_0 = 2, 3 levels. A single cross section calculation exists for the metastable 2s state, performed within the adiabatic superpromotion model [12].

The cross section $\sigma_{ion}(n_0 = 2)$ was determined by scaling the corresponding cross section for He²⁺ impact from Ref. [11] according to scaling relation [22,23]

$$\frac{\sigma_{ion}(q,n_0)}{qn_0^4} = f(n_0^2 E/q), \tag{16}$$

The cross section $\sigma_{ion}(n_0 = 3)$ at high energies was determined from the scaled BBA formula (see [11]), and at low energies (below $\sim 20 \text{keV/amu}$) by using the ratio

$$\frac{\sigma_{ion}(Li^{3+}; n_0 = 3)}{\sigma_{ion}(Li^{3+}; n_0 = 2)} = \frac{\sigma_{ion}(He^{2+}; n_0 = 3)}{\sigma_{ion}(He^{2+}; n_0 = 2)}$$
(17)

in which $\sigma_{ion}(He^{2+}; n_0 = 2,3)$ were taken from Ref. [11].

The cross sections $\sigma_{ion}(n_0=2,3)$ derived in this way have been fitted to an expression of the form given by Eq. (15). The corresponding fitting coefficients a_i are given in Table 7. The uncertainty of $\sigma_{ion}(n_0=2,3)$ cross sections is estimated to be 30-40% for E\leq10 keV/amu, 20-30% for $10 \le E(\text{keV/amu}) \le 100$, and $\sim 20\%$ for $E \ge 100 \text{ keV/amu}$.

3.2.2. Ionization from $n_0 \ge 4$ levels

The cross section $\sigma_{ion}(n_0 \ge 4)$ can be derived by using the scaling relation

$$\sigma_{ion}(n_0; E_{n_0}) = \left(\frac{n_0}{3}\right)^4 \sigma_{ion}(n_0 = 3; E), \quad E_{n_0} = \left(\frac{3}{n_0}\right)^2 E$$
 (18)

where $\sigma_{ion}(n_0 = 3; E)$ is the ionization cross section for $n_0 = 3$ level given in the preceding sub-section. The scaling relation (18) should be fairly accurate ($\sim 10\%$ uncertainty) for $E n_0 \ge 100 \text{ keV/amu}$, but it becomes less accurate at lower energies.

4. Electron Capture

4.1. Electron capture from H(1s)

Electron capture in Li³⁺ + H(1s) collisions has been subject to many both

theoretical and experimental studies. The motivation for this is the fact that at low and intermediate collision energies the electron capture process populates dominantly the excited states of $\text{Li}^{2+}(n\ell)$ ion, the radiation from which can be used for diagnostic purposes (see, e.g., [1,2]). As a result of these studies, both the total and state-selective (for a given $n\ell$), cross sections for this process are known to a high accuracy in the energy range from $\sim 100 \text{ eV/amu}$ to 10 MeV/amu.

4.1.1. Total cross section

There have been two experimental measurements of the total cross section of $\text{Li}^{3+} + \text{H}(1\text{s})$ charge exchange reaction covering the energy ranges $15 \cdot 250$ keV/amu[14] and $\sim 1 \cdot 6$ keV/amu [24]. Theoretical calculations are numerous; they employ different types of methods with varying accuracy and taken together, cover the energy range from 80 eV/amu to 10 MeV/amu. The most accurate of these calculations agree with each other, as well as with the experimental data (except for the lower measured energies of Ref. [24]). The total charge exchange (CX) cross section is sum over the electron capture cross sections to specific (n, ℓ) final states, $\sigma_{CX}^{(n,\ell)}$, of product ion $\text{Li}^{2+}(n\ell)$,

$$\sigma_{CX}^{tot} = \sum_{n,\ell} \sigma_{CX}^{(n,\ell)} = \sum_{n} \sigma_{CX}^{(n)}$$

$$\tag{19}$$

where

$$\sigma_{CX}^{(n)} = \sum_{\ell} \sigma_{CX}^{(n,\ell)} \tag{20}$$

is the electron capture cross section to all ℓ -sub-states within a given n-shell. Since in most applications (e.g. CX recombination spectroscopy, CXRS) $\sigma_{CX}^{(n,\ell)}$ are of main interest, we present in the next sub-section the best available theoretical information for $\sigma_{CX}^{(n,\ell)}$. The total cross section σ_{CX}^{tot} is then obtained from Eq. (19).

4.1.2. $\sigma_{CX}^{(n,\ell)}$ and $\sigma_{CX}^{(n)}$ partial cross sections

While experimental measurements of state-selective electron capture cross sections $\sigma_{CX}^{(n,\ell)}$ in reaction (3) have not yet been reported, there have been many theoretical calculations of these cross sections, several of them having high accuracy ($\leq 10\%$ uncertainty) in the energy region of their applicability.

In the low energy region (E \leq 20keV/amu), accurate calculations of $\sigma_{CX}^{(n,\ell)}$ have been performed by the advanced adiabatic (superpromotion) method [5] and by the MO-CC method employing large number of MOs [25,26]. The calculations in Ref. [26] go down to energies of ~80 eV/amu. In the intermediate-to-high energy region (E = 15-200 keV/amu), there exist a number of large-base AO-CC calculations of $\sigma_{CX}^{(n,\ell)}$ [6,27,27a,28], some of them employing also a large number of pseudo-states [27,27a,28]. In this energy range (but above 37.5 keV/amu) there exist also extensive CTMC calculations [29], some of them being performed (for the n=6 states) in the course of the present work. For the capture to 2s state, a cross section calculation has been performed in Ref. [30] by a second-order type method. In the high-energy region (E \geq 100 keV/amu), accurate $\sigma_{CX}^{(n,\ell)}$ cross sections for the reaction (3) have been performed within the second-order continuum distorted wave (CDW) approximation [31] up to 10 MeV/amu, and by using other less sophisticated methods [32-34]. For $\sigma_{CX}^{(n)}$, CDW data are also given in [35]. The most accurate of these calculations are those of Refs. [25,26] for low energies, Refs. [27,27a,28] for intermediate-to-high energies, and Ref. [31] for the high-energy region. In the energy range of applicability of the employed method, all these calculations have an uncertainty of about 10%, or smaller, at least for the dominant channels. For the electron capture (n,ℓ) channels having small cross sections, the accuracy is not expected to be so high. The cross section results from these references have been taken as a basis for constructing the $\sigma_{CX}^{(n,\ell)}$ cross sections in the energy region from \sim 80 eV/amu to 10 MeV/amu. Since the energy ranges in which the above calculations been made have a significant overlap (mainly due to validity of AO-CC method with inclusion of pseudo-states in the broad region from -10-200 keV/amu), a smooth connection of the results from different energy regions was possible.

The $\sigma_{CX}^{(n)}$ partial cross sections show the following properties. In the energy region below $\sim 2 \text{ keV/amu}$, $\sigma_{CX}^{(n=3)}$ cross section is dominant, while for E > 2

keV/amu, the cross sections $\sigma_{CX}^{(n=2)}$ and $\sigma_{CX}^{(n=3)}$ have comparable magnitudes and are the largest ones. The cross sections $\sigma_{CX}^{(n)}$ for n=4,5,6,7··· rapidly decrease with increasing n (particularly at energies below ~ 50 keV/amu; for energies above ~ 200 keV/amu this decrease follows the n^{-3} Oppenheimer's rule). The $\sigma_{CX}^{(n=1)}$ cross section is very small: in the region of its maximum at E ≈ 100 keV/amu it is of order of magnitude of 10^{-18} cm² and is about an order of magnitude smaller than $\sigma_{CX}^{(n=7)}$, the higher n for which state-selective calculations have been performed.

The $\sigma_{CX}^{(n,\ell)}$ cross sections show a more complex dependence on parameter ℓ for a given n and energy. At very low energies ($\leq 1 \text{ keV/amu}$), the maximum populated ℓ -sub-level for a given n is $\ell_m = n-1$. For n=2,3, the dominance of $\sigma_{CX}^{(n,n-1)}$ persists even up to $E \cong 300 \text{keV/amu}$. With increasing both the energy and n, however, the maximum of ℓ -distribution of captured electrons (i.e. the ratio $\sigma_{CX}^{(n,\ell)}/\sigma_{CX}^{(n)}$) shifts towards lower ℓ -values (for instance, for n=4 and $1 \leq 100$, $1 \leq 100$,

The constructed cross sections $\sigma_{CX}^{(n,\ell)}$ on the basis of the data from Refs. [25·28,31] (considered to be the best) have been fitted to the following types of analytic expressions, the different form of which reflects the different energy behavior of the cross sections

$$\sigma_{CX}^{(n,\ell)} = \frac{a_1 \exp\left[-\left(a_2/E\right)^{a_3}\right]}{1 + \left(E/a_4\right)^{a_5} + \left(E/a_4\right)^{a_7}} \quad \left(\times 10^{-16} cm^2\right)$$
(21)

$$\sigma_{CX}^{(n,\ell)} = \frac{a_i \exp\left[-\left(a_2/E\right)^{a_3}\right]}{1 + \left(E/a_4\right)^{a_5} + \left(E/a_6\right)^{a_7} + \left(E/a_8\right)^{a_9}} \quad \left(\times 10^{-16} \, cm^2\right)$$
(22)

$$\sigma_{CX}^{(n,\ell)} = \frac{a_1 \exp\left[-\left(a_2/E\right)^{a_3}\right]}{1 + \left(E/a_4\right)^{a_5} + \left(E/a_6\right)^{a_7} + \left(E/a_8\right)^{a_9}} + \frac{a_{10} \exp\left[-\left(a_{11}/E\right)^{a_{12}}\right]}{1 + \left(E/a_{13}\right)^{a_{14}}} \left(\times 10^{-16} cm^2\right)$$
(23)

$$\sigma_{CX}^{(n,\ell)} = \frac{a_1 \exp\left[-\left(a_2/E\right)^{a_3}\right]}{1 + \left(E/a_4\right)^{a_5} + \left(E/a_6\right)^{a_7} + \left(E/a_8\right)^{a_9}} + \frac{a_{10} \exp\left[-\left(a_{11}/E\right)^{a_{12}}\right]}{1 + \left(E/a_{13}\right)^{a_{14}} + \left(E/a_{15}\right)^{a_{16}}} \quad \left(\times 10^{-16} \, cm^2\right) \tag{24}$$

where collision energy E is expressed in eV/amu, and a_i (i=1-16) are fitting parameters. The values of fitting parameters a_i for $\sigma_{CX}^{(n,\ell)}$ cross sections for n=1 and n=2, and n=3-6 are given in Table 8, and Table 9-12, respectively. The number of parameters in these tables indicates the used analytic fit expression from those given by Eqs. (21)-(24). We note that $\sigma_{CX}^{(n,\ell)}$ for $n\ell=6h$ is small, and corresponding values of parameters a_i are not shown in Table 12.

The shell-selective partial cross sections $\sigma_{CX}^{(n)}$, obtained by summing $\sigma_{CX}^{(n,\ell)}$ over ℓ , can also be represented by an analytic fit expression in the same form as Eq. (21-24). The corresponding values of fitting parameters a_i for n = 2-7 are given in Table 13.

For all the above electron capture processes, the available theoretical results and the fitting line for the preferred ("recommended") cross section are displayed in figures 3-28.

The extension of $\sigma_{CX}^{(n)}$ and $\sigma_{CX}^{(n,\ell)}$ data to higher $n \geq 8$ cannot be done because of absence of appropriate scalings. Only in the high-energy region (E $\geq 200 \text{ keV/amu}$), the Oppenheimer's n^{-3} rule for $\sigma_{CX}^{(n)}$ is known to be valid [36].

4.2. Capture from $H(n_0 \ge 2)$ levels

Electron capture in H(2s)+ Li³⁺ collisions has been studied at low collision energies by the Landau-Zener [37] and superpromotion [5,38] models, and at intermediate to high energies by the AO-CC [39] and other less accurate models [40]. The total electron capture cross section shows a broad maximum in the energy region 0.5·5 keV/amu of about $10^{-14} cm^2$. For energies below $\sim 0.5 \cdot 5$ keV/amu, dominantly populated final n-shells are n=5 and n=6 ($\sigma_{CX}^{(n=6)}$ is, however, by a factor of two smaller than $\sigma_{CX}^{(n=5)}$), while for $E \geq 0.5$ keV/amu, n=4 shell is dominantly populated. Within n=4 shell, the 4f state is dominantly populated up to $E \approx 5$ keV/amu and then its population becomes comparable with that of 4d. Within n=5, dominantly populated state is 5f for energies below ~ 10 keV/amu.

The state-selective electron capture cross sections for H(2p)+Li³⁺ collisions have been calculated only by the AO-CC method [39]. The cross section magnitude and relative importance of individual final (capture) channels are

similar as in the case of H(2s)+Li³⁺ case.

For $n_0 \ge 2$, total electron capture cross section calculations are available (for selected n_0 between $n_0 = 2$ and $n_0 = 20$) from the CTMC calculations of Ref. [41] (for q = 2) and Ref. [42] (q = 5,10). These data, together with the experimental data on $H(n_0) + He^{2+}$ collision system ($n_0 = 8-24$) [43], have been represented in Ref. [44] in a compact, scaled form (within an uncertainty of 10-30%)

$$\widetilde{\sigma}_{CX}^{tot}(q) = \frac{\sigma_{CX}^{tot}(n_0)}{qn_0^4} = 7.04 \frac{a}{\widetilde{E}^{3.5}(1 + b\widetilde{E}^2)} \left\{ 1 - \exp\left[-\frac{2\widetilde{E}^{3.5}(1 + b\widetilde{E}^2)}{3a} \right] \right\} \quad (\times 10^{-16} cm^2) \quad (25)$$

where $\widetilde{E} = E n_0^2 / q^{0.5}$ (keV/amu), $a = 1.507 \times 10^5$, $b = 1.974 \times 10^{-5}$.

Below $\widetilde{E} \sim 10 \text{ keV/amu}$, the energy dependence of σ_{CX}^{tot} is very weak, and for $\widetilde{E} \rightarrow 0$, $\widetilde{\sigma}_{CX}^{tot} = 1.06 \times 10^{-15} cm^2$. The experimental data of Ref. [43] for $H(n_0 \ge 8) + He^{2+}$, represented by Eq. (25) with an accuracy of $\sim 20\%$, extend down to $\widetilde{E} \approx 10 \text{ keV/amu}$. For the partial cross sections $\sigma_{CX}^{(n)}(n_0)$ and $\sigma_{CX}^{(n,\ell)}(n_0)$ with $n_0 \ge 3$ neither direct calculations nor scaling relationships are available. From the general theory of charge exchange reactions of one-electron systems with multiply charged ions [36] it follows that dominantly populated n-shell of $\text{Li}^{2+}(n)$ at energies below $\sim 1/n_0 \text{ keV/amu}$ should be [36]

$$n_m \equiv 2^{1/4} n_0 q^{3/4} \equiv 2.71 n_0 \tag{26}$$

5. Conclusions

In the present report we have compiled and critically assessed the available cross section information on collision processes of Li³⁺ ions with atomic hydrogen. The processes include excitation, ionization and charge exchange from the ground and excited hydrogen levels. The energy range considered extends from $\sim 100 \text{eV/amu}$ to the MeV/amu region.

The cross section information on these processes found in literature pertains mostly to H in its ground state. For the processes involving excited initial states it is either sparse (the case of charge exchange) or completely missing (for excitation and ionization). Therefore, certain well established cross section scaling relationships had to be employed to generate the missing cross section information.

In the case of charge exchange process from H(1s), capture to specific states of Li^{2+} ($n\ell$) ion has been also considered in view of its importance for plasma charge exchange recombination diagnostics.

The critically assessed cross sections for above processes have been represented by relatively simple analytic fit functions. These fit functions have been chosen to have correct physical behavior at both low and high energies and can ensure reliable extrapolations. The error of the analytic fits to the cross sections is well within the uncertainty of original data.

Although the present database can usefully serve the needs for Li³⁺ -based plasma diagnostics, its completion with the presently missing cross sections (especially for electron capture from excited $H(n \ge 2)$ levels) is highly desirable.

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Table 1. Values of parameters a_i in Eq. (5) for excitation of n=2,3,4 levels

a_i	n=2	n=3	n=4
a_{l}	4.84520(+2)*	5.04618(+1)	1.59325(+1)
a_2	6.69602(+1)	8.21366(+1)	7.94188(+1)
a_3	2.78499(-2)	4.50894(-1)	8.82429(-1)
a_4	1.01627(-5)	7.06132(-6)	6.52302(-6)
a_{5}	4.84603(-3)	5.16969(-1)	4.46519(-1)
a_6	-3.19529(+0)	-4.17842(†0)	-4.24975(†0)
a_7	3.36324(-2)	1.67102(-1)	4.12196(-2)
a_8	2.64997(+1)	2.87789(+1)	2.28376(+1)
a_9	9.56579(-4)	1.59714(-4)	1.58302(-8)
a_{10}	3.84978(+0)	4.77936(†0)	8.74347(+0)

 $[\]star \overline{a(\pm x)} = \overline{a \times 10^{\pm x}}$

Table 2. Values of fitting parameters b_i in Eq. (6) for excitation of n = 5,6 levels

b_{i}	n=5	n=6
b_1	1.1204(+1)*	5.7394(+0)
b_2	9.6873(+1)	1.1152(+2)
b_3	7.0587(-2)	1.3088(-1)
b_4	7.2281(-5)	8.8253(-5)
b_5	1.0354(-1)	8.5857(-2)
b_6	-2.4161(+0)	-2.2905(+0)
b_7	1.9046(+0)	1.9145(+0)
b_8	-6.3726(+0)	-5.5986(+0)

 $[*]a(\pm x) = a \times 10^{\pm x}$

Table 3. Values of parameters c_i in Eq. (8) for the excitation transition of $n_0 = 2 \rightarrow n = 3,4,5$

c_i	n=2	n=3	n=4
c_1	3.95830(+3)*	5.26643(†2)	1.73597(+2)
c_2	4.67919(+1)	5.18702(+1)	5.69191(+1)
c_3	1.72288(-1)	6.21216(-1)	9.68539(-1)
C 4	1.42039(-3)	7.86601(-4)	9.00919(-4)
c_{5}	1.08533(-1)	1.17296(-1)	1.05702(-1)
c_6	-1.22181(+0)	-1.56016(+0)	-1.47750(+0)

^{*} $a(\pm x) = a \times 10^{\pm x}$

Table 4. Values of coefficients c_{2-n} in Eq. (9) for excitation transition of $2 \rightarrow n(n = 6,7,8,9,10)$

n	n=6	n=7	n=8	n=9	n = 10
C_{2-n}	0.4610	0.2475	0.1465	0.0920	0.0605

Table 5. Values of fitting parameters c_i in Eq. (8) for $n_0 = 3 \rightarrow n = 4.5.6$ excitation transitions

	<u>. V</u>		
c_{i}	n=4	n=5	n=6
c_1	1.16987(+4)*	1.78730(+3)	5.95331(+2)
c_2	3.39570(+1)	3.32880(+1)	3.45210(+1)
c_3	8.74514(-1)	9.69935(-1)	1.44722(+0)
C 4	5.66708(-3)	5.92691(-3)	6.31140(-3)
c_{5}	1.73728(-1)	1.80588(-1)	1.78202(-1)
c 6	-1.27220(+0)	-1.28940(+0)	-1.28810(+0)

 $[*] a(\pm x) = a \times 10^{\pm x}$

Table 6. Values of coefficients c_{3-n} in Eq. (11) for excitation transitions $3 \rightarrow n(n=7-10)$

n	n=7	n=8	n=9	n=10
C_{3-n}	0.4670	0.2545	0.1540	0.1000

Table 7. Values of fitting parameters a_i in Eq. (15) for ionization cross section of $H(n_0)$, n = 1,2,3 by Li^{3+} impact

a_{i}	$n_0 = 1$	$n_0 = 2$	$n_0 = 3$
a_1	8.30194(+1)*	1.66484(†2)	7.53658(+2)
a_2	1.15973(+2)	5.19714(+1)	1.92724(+1)
a_3	2.77580(+7)	4.52808(+21)	1.83610(+9)
a_4	4.59930(-6)	2.16126(-3)	4.93943(-2)
a_5	1.91247(-2)	4.74902(-2)	1.36404(-1)
a_6	-3.57909(+0)	-2.41416(+0)	-1.16090(+0)
a_7	6.29546(-2)	1.60611(-3)	1.30281(+0)
a_8	-1.62357(+0)	3.88616(-2)	-2.77601(+0)

 $[\]star \overline{a(\pm x)} = a \times 10^{\pm x}$

Table 8. Values of fitting parameters a_i for $\sigma_{CX}^{(n,\ell)}$ in Eq. (23) for n=1 and in Eq. (22) for n=2

	III Eq. (20	7101 / T dild li	$\frac{1}{n}$ Eq. (22) for n -
a_{i}	1s	2s	2p
a_1	8.5833(-1)	1.0795(+3)	1.8660(+3)
a_2	1.0919(+6)	1.2027(+6)	1.2061(+6)
a_3	4.6504(-1)	2.9973(-1)	2.9785(-1)
a_4	2.9509(+4)	4.1258(+3)	1.5605(+3)
a_5	-2.1907(+0)	1.7088(+0)	8.6085(-1)
a_6	9.5881(+4)	9.3273(+3)	1.6528(+4)
a_{7}	3.1259(+0)	3.1257(+0)	3.4454(+0)
a_8	2.4012(+5)	4.9259(+4)	5.3094(+4)
a_9	5.6032(+0)	5.6066(+0)	6.1615(+0)
a_{10}	1.6874(+1)		
a_{11}	1.3618(+4)		
a ₁₂	6.7598(-1)		
a_{13}	2.1373(+1)		
a_{14}	3.5534(+0)		

 $[*] a(\pm x) = a \times 10^{\pm x}$

Table 9. Values of fitting parameters a_i for $\sigma_{CX}^{(n,\ell)}$ in Eq. (22) for 3s and 3p, and in Eq. (23) 3d

a_i	3s	3p	3d
a_{i}	1.2501(+1)	2.3666(+1)	2.2209(+1)
a_2	2.9346(+6)	2.0288(+6)	1.8692(+4)
a_3	1.5702(-1)	1.5563(-1)	5.5415(-1)
a_4	5.3156(+4)	5.9278(+4)	5.7957(+4)
a_5	4.3742(+0)	4.5574(+0)	6.0668(+0)
a_6	1.7553(+4)	4.1115(+4)	2.9885(+4)
a_7	2.1353(+0)	2.5256(+0)	3.4442(+0)
a_8	1.1964(+5)	9.7507(+4)	9.7442(+4)
a_9	5.9502(+0)	6.1913(+0)	5.9651(+0)
a_{10}			6.0765(+0)
a_{11}			2.1264(+3)
a_{12}			2.7457(-1)
a_{13}			1.0146(+3)
a_{14}			3.2956(+0)

^{*} $a(\pm x) = a \times 10^{\pm x}$

Table 10. Values of fitting parameters a_i in Eq. (24) for $\sigma_{CX}^{(n,\ell)}$ for n=4

a_i	4s	4p	4d	4f
a_1	1.9807(+0)	5.8172(+0)	6.8352(+0)	2.6988(+0)
a_2	3.0693(+5)	3.0879(+5)	3.0879(+5)	3.0879(+5)
a_3	3.1672(-1)	3.3179(-1)	3.2622(-1)	3.1213(-1)
a_4	7.8964(+4)	7.7848(+4)	7.7853(+4)	7.7831(+4)
a_5	3.3043(+0)	4.2739(+0)	3.0443(+0)	5.8111(+0)
a_6	3.9049(+4)	6.4526(+4)	6.4635(+4)	6.4459(+4)
a_7	2.8297(+0)	4.0548(+0)	5.5804(+0)	6.0667(+0)
a_8	8.7775(+4)	9.5977(+4)	9.5968(+4)	9.5973(+4)
a_9	5.4039(+0)	6.1357(+0)	6.4213(+0)	1.9245(+0)
a_{10}	1.7856(+2)	2.7492(-1)	9.8665(-1)	5.2029(-1)
a_{11}	5.0490(+3)	4.8559(+3)	4.8536(+3)	4.8690(+3)
a_{12}	3.0792(+0)	1.2215(+0)	1.3517(+0)	7.2433(-1)
a_{13}	1.0748(+3)	1.7250(+3)	1.7210(+4)	1.6908(+3)
· a ₁₄	1.3623(+1)	1.1163(+0)	1.7927(+0)	6.5445(-1)
a_{15}	7.5998(+3)	7.5993(+3)	7.5989(+3)	7.6791(+3)
a_{16}	5.9852(+0)	4.2605(+1)	5.2865(+1)	4.0681(+0)

^{*} $a(\pm x) = a \times 10^{\pm x}$

Table 11. Values of fitting parameters a_i in Eq. (24) for $\sigma_{CX}^{(n,\ell)}$ for n=5

a_i	5s	5p	5d	5f	5g
a_1	2.9589(+4)	2.9610(+4)	2.9560(+4)	2.9551(+4)	2.9588(+4)
a_2	8.6902(+5)	8.6902(+5)	8.6902(+5)	8.6902(+5)	8.6902(+5)
a_3	4.7454(-1)	5.1531(-1)	4.3330(-1)	4.2289(-1)	4.5340(-1)
a_4	6.1427(+3)	6.1425(+3)	6.1423(+3)	6.1423(+3)	6.1427(+3)
a_{5}	5.8954(-1)	1.1564(+0)	8.5495(-1)	1.5005(+0)	6.5271(-2)
a_6	6.4719(+2)	5.9453(+2)	7.4504(+1)	1.0527(+2)	8.0194(+2)
a_7	2.3567(+0)	1.4956(+0)	1.2092(+0)	1.5162(+0)	2.6179(+0)
a_8	1.9494(+4)	1.9694(+4)	1.9219(+4)	1.9136(+4)	1.9467(+4)
a_9	5.8919(+0)	. 6.0080(+0)	6.2439(+0)	6.7015(+0)	7.5035(+0)
'a ₁₀	8.8685(-2)	9.2030(-2)	4.3839(-2)	3.1062(-2)	7.2181(-2)
a_{11}	4.1267(+4)	4.1269(+4)	4.1287(+4)	4.1283(+4)	4.1271(+4)
a ₁₂	4.0393(-1)	3.3759(-1)	4.1175(-1)	3.9622(-1)	4.0936(-1)
<i>a</i> ₁₃	1.1251(+3)	1.3872(+3)	1.1241(+4)	1.1240(+4)	1.1255(+4)
a ₁₄	-5.5447(-1)	-1.9538(+0)	2.0048(+0)	2.2867(+0)	2.3253(+0)
a ₁₅	6.1107(+4)	6.1108(+4)	6.1148(+4)	6.1149(+4)	6.1124(+4)
a_{16}	4.7226(+0)	5.8212(+0)	5.2734(+0)	5.0499(+0)	5.4616(+0)

 $[*] a(\pm x) = a \times 10^{\pm x}$

Table 12. Values of fitting parameters a_i in Eq. (21) for $\sigma_{CX}^{(n,\ell)}$ for n=6 ($\sigma_{CX}^{(n,\ell)}$ for $n\ell=6h$ is small and not shown here)

		** ***********************************			
a_i	6s	6p	6d	6f	6g
a_1	2.8852(+5)*	4.1604(+5)	2.5759(+4)	2.5605(+4)	2.2209(+4)
a_2	7.6028(+5)	1.1919(+6)	1.1456(+5)	4.4935(+5)	2.6385(+5)
a_3	6.4890(-1)	5.8398(-1)	9.5729(-1)	7.4489(-1)	8.3042(-1)
a_4	5.0200(+3)	4.5375(+3)	1.6186(+2)	9.4048(+3)	8.9336(+3)
a_{5}	4.3527(+0)	3.2608(+0)	1.6786(+0)	4.0746(+0)	5.7968(+0)
a_6	1.1099(+4)	1.6024(+4)	1.3463(+4)	1.2281(+4)	1.6229(+4)
a_7	5.7395(+0)	6.3637(+0)	5.9743(+0)	5.9934(+0)	5.8128(+0)

 $[*] a(\pm x) = a \times 10^{\pm x}$

Table 13. Values of fitting parameters a_i in Eq. (21-24) but for n-selective cross sections $\sigma_{CX}^{(n,\lambda)}$, for n=2-7

a_{i}	2	3	4	5	6	7
$a_{\rm l}$	3.2285(+3)	4.5013(+5)	2.7945(+1)	3.9605(+4)	1.0645(+0)	6.8054(+4)
a_2	1.2032(+6)	6.0007(-3)	3.0879(+5)	8.6902(+5)	5.9413(+4)	8.4908(+5)
a_3	2.9923(-1)	-1.5258(-1)	3.2323(-1)	4.6249(-1)	5.1555(-1)	4.8405(-1)
a_4	2.4583(+3)	1.3138(+4)	7.7827(+4)	6.1422(+3)	7.9956(+4)	3.3952(+2)
a_5	1.2528(+0)	-3.4971(+0)	3.9735(+0)	1.2053(+0)	4.4936(+0)	1.7087(+0)
a_6	1.8499(+4)	7.2544(+4)	6.4553(+4)	4.6778(+2)	1.0236(+5)	1.6477(+4)
a_7	3.9143(+0)	2.5157(+0)	5.5443(+0)	1.4458(+0)	5.7539(+0)	5.8670(+0)
a_8	4.8151(+4)	2.1012(+4)	9.5977(+4)	1.9698(+4)		
a_9	5.7723(+0)	-1.3108(+0)	1.6934(-1)	5.7783(+0)		
a_{10}		1.4577(+1)	3.8125(+0)	3.0066(-1)		
a_{11}	-	1.8457(+5)	4.8562(+3)	4.1268(+4)		
a_{12}		1.3262(-1)	1.1121(+0)	3.2486(-1)		
<i>a</i> ₁₃		4.4535(+3)	1.7255(+3)	1.3114(+3)		<u></u>
a ₁₄		3.5393(+0)	2.0477(+0)	-1.9767(+0)		·
a_{15}			7.6001(+3)	6.1110(+4)		·
<i>a</i> ₁₆			1.8493(+1)	5.6313(+0)		

 $[*]a(\pm x) = a \times 10^{\pm x}$

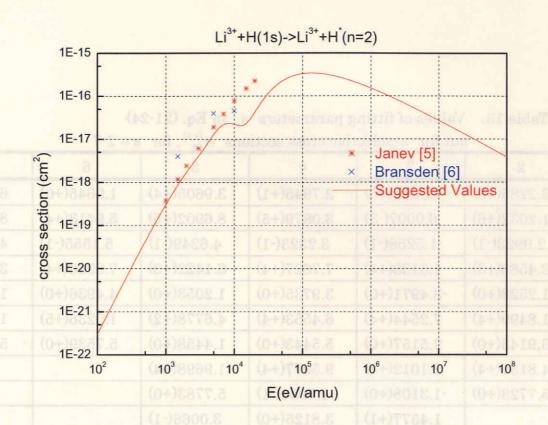


Figure 1. Preferred (full line; see text) cross section for excitation of H(n=2) by Li³⁺ colliding with H(1s).

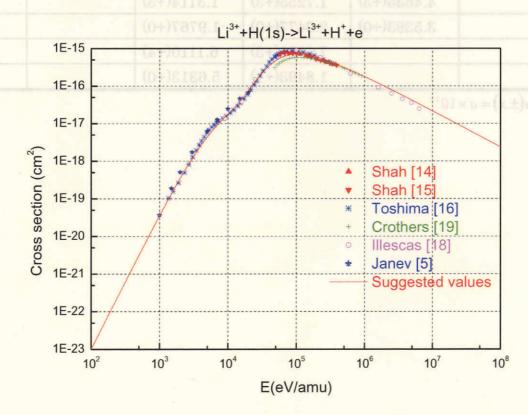


Figure 2. Fit (full line) of preferred ionization cross section of H(1s) by Li³⁺ impact.

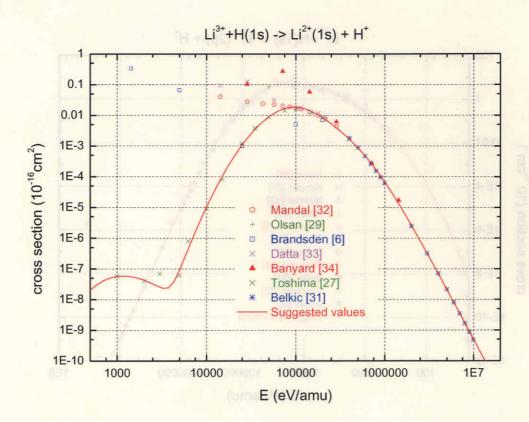


Figure 3. Fit (full line) of preferred electron capture cross section to Li²⁺(1s) in collisions of Li³⁺ with H(1s).

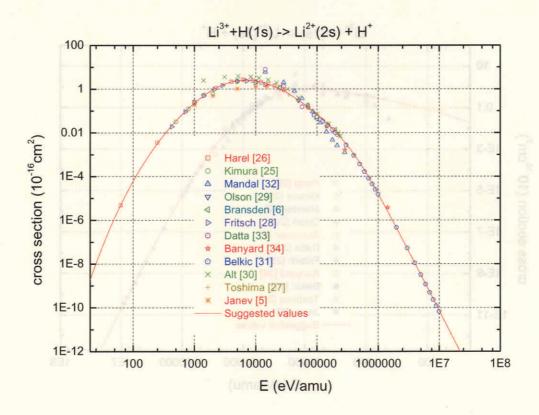


Figure 4. Same as in Fig.3, but for capture to Li²⁺(2s).

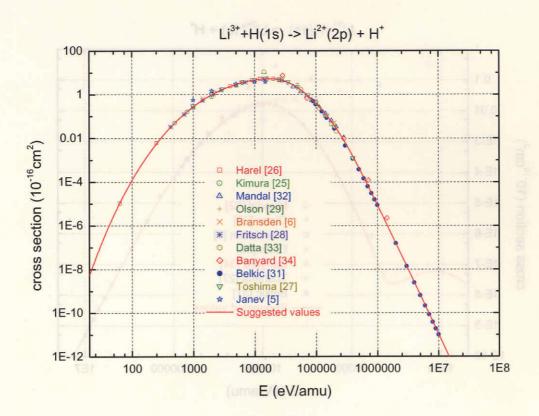


Figure 5. Same as in Fig.3, but for capture to Li²⁺(2p).

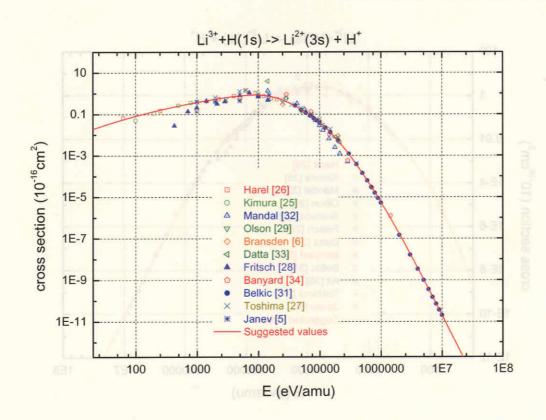


Figure 6. Same as in Fig.3, but for capture to Li²⁺(3s).

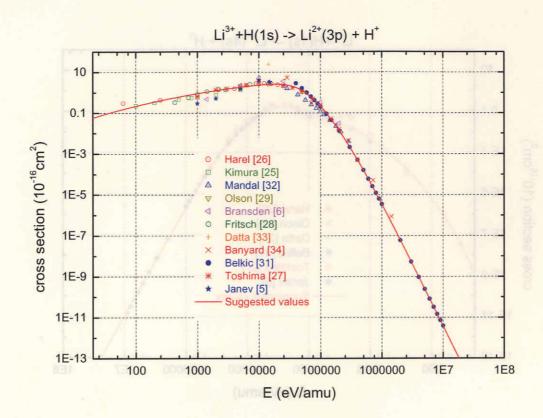


Figure 7. Same as in Fig.3, but for capture to Li²⁺(3p).

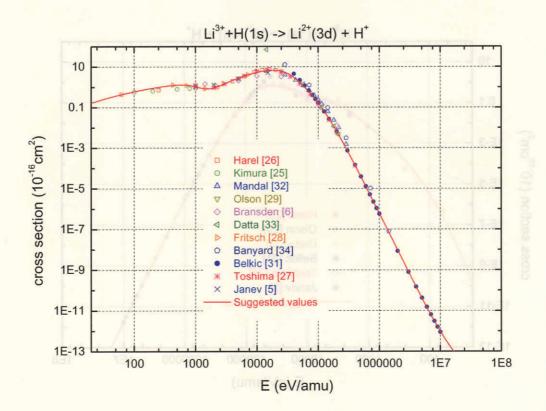


Figure 8. Same as in Fig.3, but for capture to Li²⁺(3d).

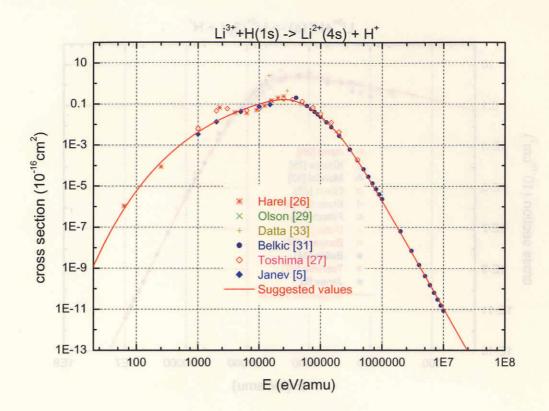


Figure 9. Same as in Fig.3, but for capture to Li²⁺(4s).

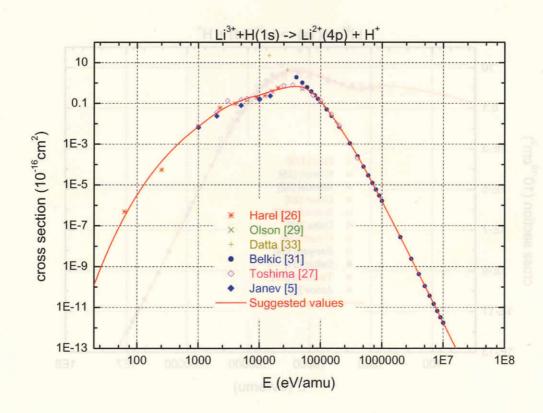


Figure 10. Same as in Fig.3, but for capture to Li²⁺(4p)

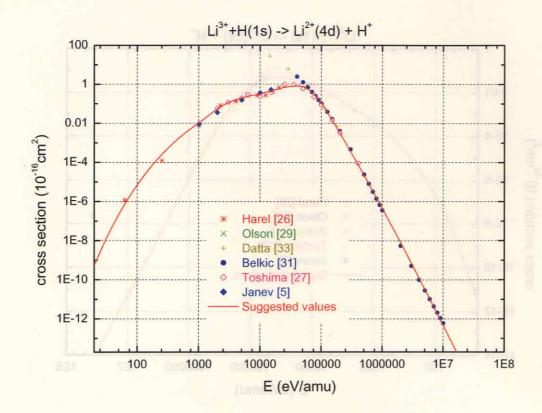


Figure 11. Same as in Fig. 3, but for capture to Li²⁺(4d).

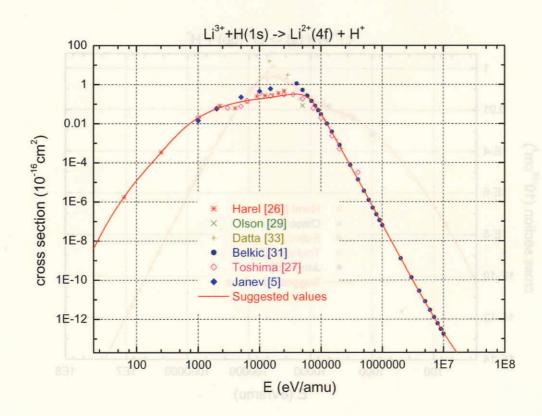


Figure 12. Same as in Fig.3, but for capture to Li²⁺(4f).

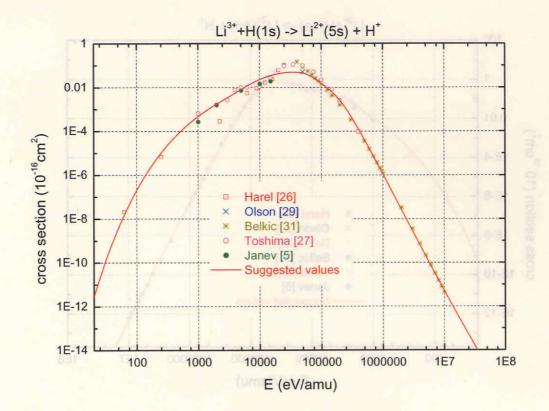


Figure 13. Same as in Fig.3, but for capture to Li²⁺(5s).

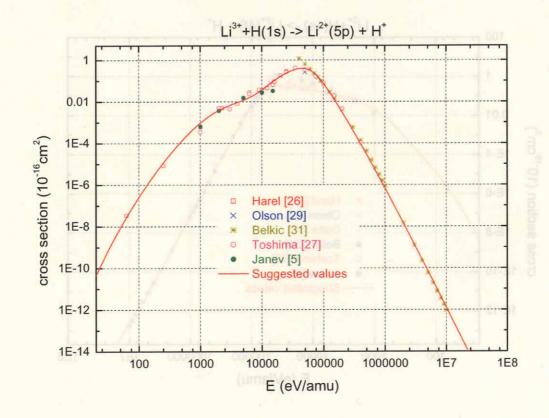


Figure 14. Same as in Fig.3, but for capture to Li²⁺(5p).

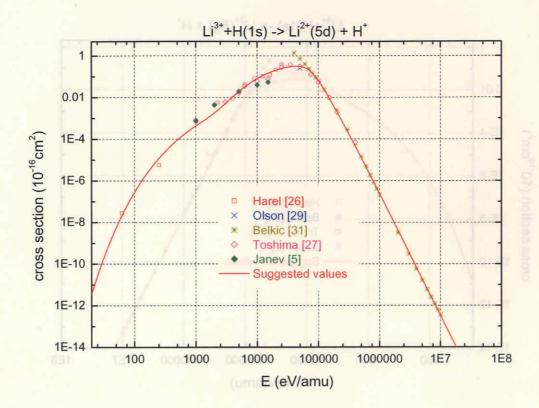


Figure 15. Same as in Fig.3, but for capture to Li²⁺(5d).

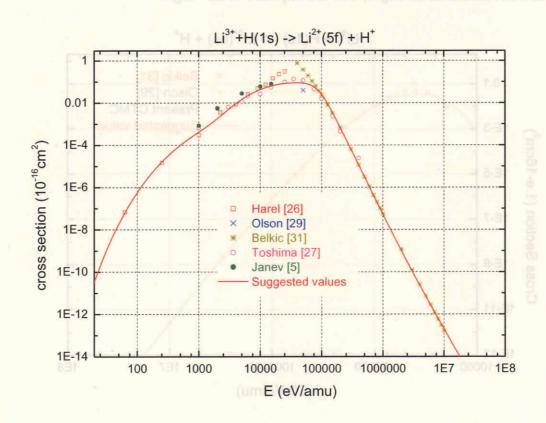


Figure 16. Same as in Fig.3, but for capture to Li²⁺(5f).

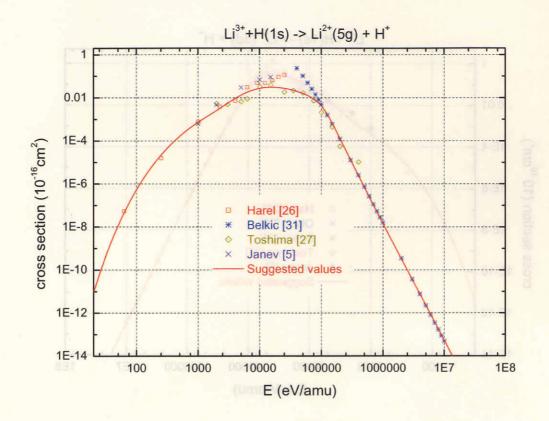


Figure 17. Same as in Fig.3, but for capture to Li²⁺(5g).

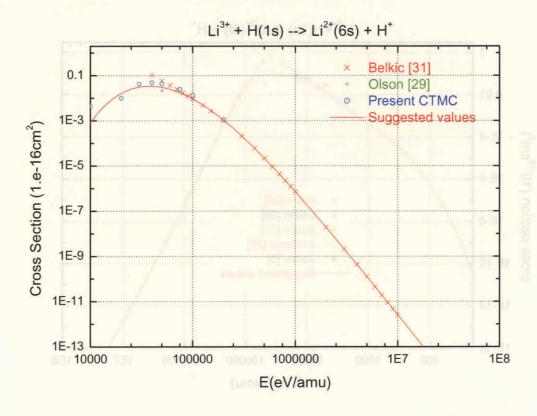


Figure 18. Same as in Fig.3, but for capture to Li²⁺(6s).

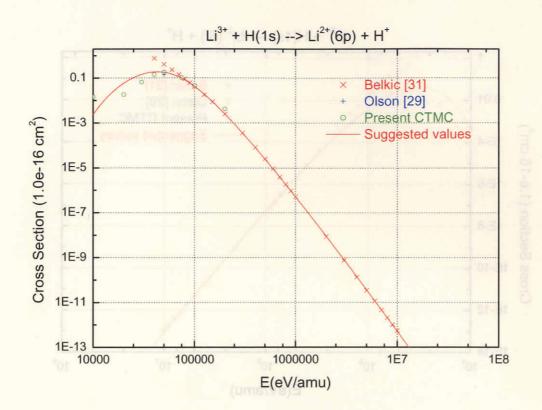


Figure 19. Same as in Fig.3, but for capture to Li²⁺(6p).

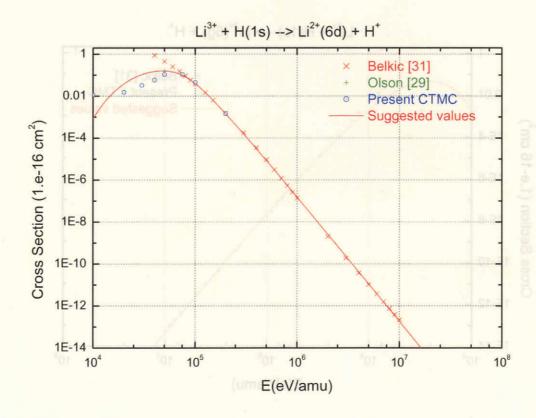


Figure 20. Same as in Fig.3, but for capture to Li²⁺(6d).

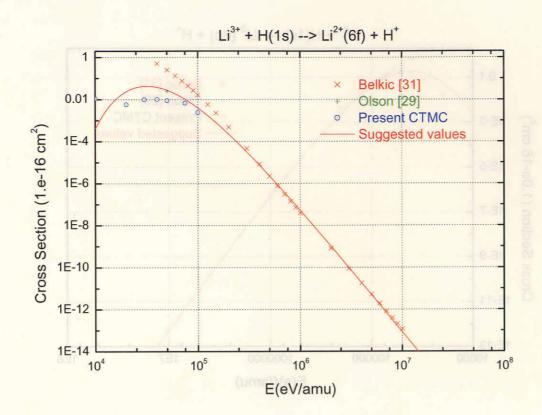


Figure 21. Same as in Fig.3, but for capture to Li²⁺(6f).

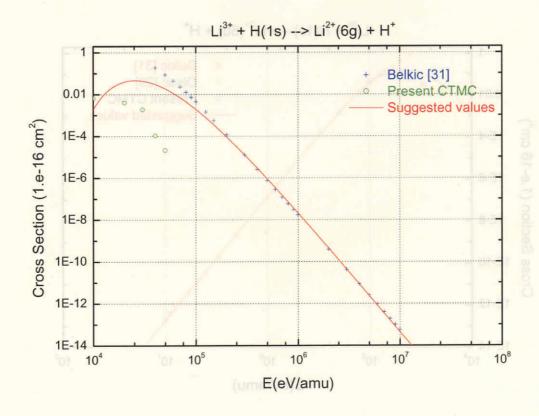


Figure 22. Same as in Fig.3, but for capture to Li²⁺(6g).

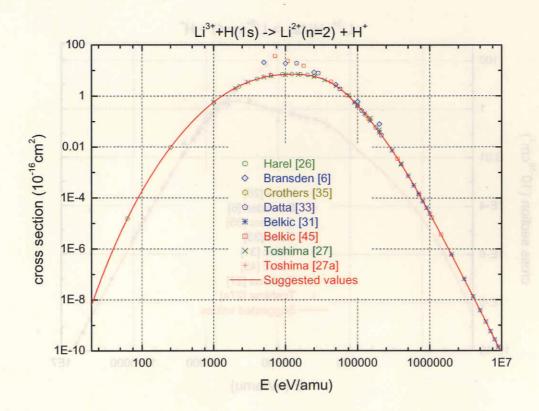


Figure 23. Fit (full line) of preferred electron capture cross section to Li²⁺(n=2) level in collisions of Li³⁺ with H(1s).

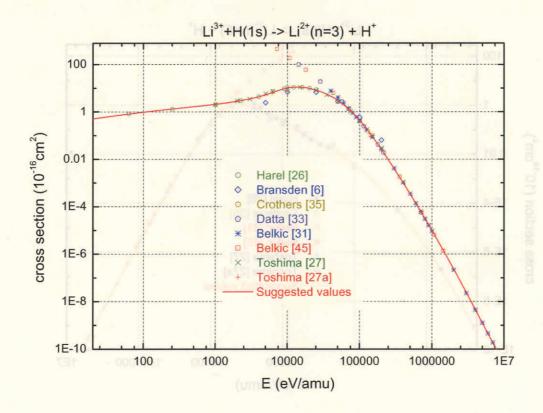


Figure 24. Same as in Fig.23, but for capture to Li²⁺(n=3).

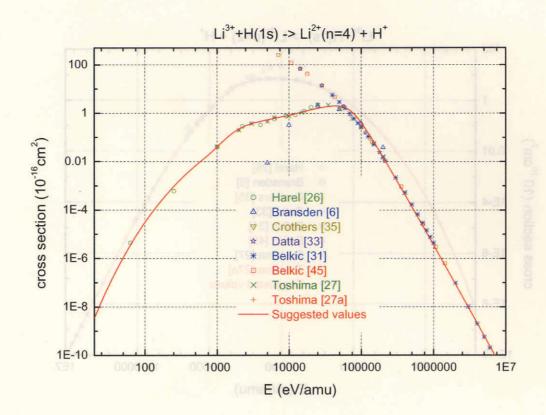


Figure 25. Same as in Fig.23, but for capture to Li²⁺(n=4).

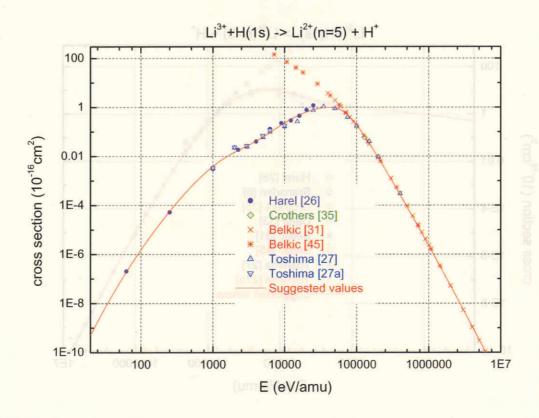


Figure 26. Same as in Fig.23, but for capture to Li²⁺(n=5).

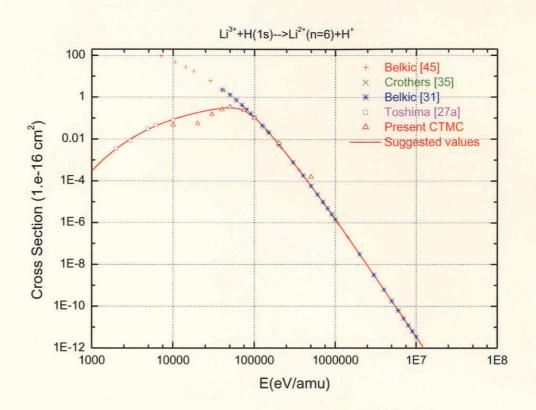


Figure 27. Same as in Fig.23, but for capture to Li²⁺(n=6).

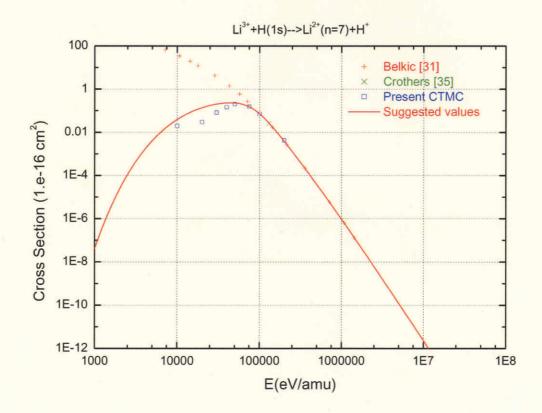


Figure 28. Same as in Fig.23, but for capture to $Li^{2+}(n=7)$.

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