

Solar fusion III: New data and theory for hydrogen-burning stars.

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In stars that lie on the main sequence in the Hertzsprung Russel diagram, like our Sun, hydrogen is fused to helium in a number of nuclear reaction chains and series, such as the proton-proton chain and the carbon-nitrogen-oxygen cycles. Precisely determined thermonuclear rates of these reactions lie at the foundation of the standard solar model.

This review, the third decadal evaluation of the nuclear physics of hydrogen-burning stars, is motivated by the great advances made in recent years by solar neutrino observatories, putting experimental knowledge of the pp chain neutrino fluxes in the few-percent precision range. The basis of the review is a one-week community meeting held in July 2022 in Berkeley, California, and many subsequent digital meetings and exchanges.

The relevant reactions of solar and stellar hydrogen burning are reviewed here, from

both theoretical and experimental perspectives. Recommendations for the state of the art of the astrophysical S-factor and its uncertainty are formulated for each of them.

Several other topics of paramount importance for the solar model are reviewed, as well: recent and future neutrino experiments, electron screening, radiative opacities, and current and upcoming experimental facilities. In addition to reaction-specific recommendations, also general recommendations are formed.

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I. Introduction

The present review summarizes the state of our understanding, in the third decade of the 21st century, of the nuclear reactions and decays taking place in solar and stellar hydrogen burning. It also addresses related issues, including solar neutrino detection, radiative opacities, electron screening of nuclear reactions, and the status of the current facilities for measuring cross sections and opacities.

As was the case for two previous decadal reviews (Adelberger *et al.*, 1998, 2011), the present review summarizes the progress made over the past decade in advancing the nuclear physics of main-sequence stars and makes recommendations for future work.

A. Purpose

Ray Davis's measurements of the flux of solar neutrinos (Davis, 2003; Davis *et al.*, 1968) showed a significant deficit with respect to the predictions of the standard solar model (SSM) (Bahcall, 1966). This discrepancy became known as the "solar neutrino problem." An intense debate ensued about possible explanations for this discrepancy, summarized in the entertaining review by Bahcall *et al.* (1988). Suggestions included plausible flaws in the SSM that might produce a cooler solar core (e.g. Schatzman *et al.* 1981; Cumming and Haxton 1996), new particle physics such as neutrino oscillations (Gribov and Pontecorvo 1969; Wolfenstein 1978), changes in the nuclear physics governing He synthesis, such as an unidentified low-energy resonance in the $^3\text{He} + ^3\text{He}$ reaction affecting extrapolations of laboratory cross sections to solar energies (Fetisov and Kopysov, 1972; Fowler, 1972), or the presence of dark matter particles modifying energy transport in the solar core (Spergel and Press, 1985). By the mid-1990s, with new results from Kamioka II/III (Hirata *et al.* 1989; Fukuda *et al.* 1996) and the SAGE/GALLEX (Abazov *et al.* 1991; Anselmann *et al.* 1992) experiments revealed a pattern of neutrino fluxes inconsistent with the expected scaling of those fluxes with the solar core temperature. Attention increasingly turned to neutrino

oscillations and other potential particle physics solutions. Concerns about the nuclear physics of the SSM also evolved, focusing more on the uncertainties that might inhibit extraction of any such "new physics."

At that time, during a meeting on the Solar Neutrino Problem hosted by the Institute for Nuclear Theory (INT), University of Washington, Seattle, a suggestion was made to convene the community working on solar nuclear reactions, in order to reach consensus on the best current values of cross sections and their uncertainties. The INT hosted the proposed workshop in February, 1997, drawing representatives from almost every experimental group active in this area, as well as many of the theorists who were engaged in solar neutrino physics. The working groups that formed during this meeting worked over the following year to evaluate past work, determining the needed cross sections and their uncertainties. The results of this evaluation, which came to be known as Solar Fusion I or SF I (Adelberger *et al.*, 1998), became the standard for use in solar modeling over the following decade.

An update of SF I was launched with the January, 2009, workshop "Solar Fusion Cross Sections II", hosted again by the INT in Seattle. Initial results from SuperKamiokande, the Sudbury Neutrino Observatory, and Borexino were then in hand, and the conversion of approximately two thirds of solar electron neutrinos into other flavors had been firmly established (Ahmad *et al.*, 2002a). Consequently, the motivation for the study was to ensure that solar model predictions would be based on the most current nuclear physics, so that meaningful uncertainties could be placed on neutrino parameters derived from solar neutrino measurements, such as the mixing angle θ_{12} . An additional source of uncertainty had arisen at the time, with the advent of three-dimensional (3D) radiative-hydrodynamic models of the Sun's atmosphere for determining element abundances from photo-absorption lines (Asplund *et al.*, 2006, 2009). This improved analysis, though, lowered the inferred metallicities, leading to tension with the Sun's interior sound speed profile determined from helioseismology (Christensen-Dalsgaard, 2002). The discrepancy between higher metallicity models that accurately reproduce the Sun's interior sound speed and lower metallicity models based on the most current treatment of the photosphere was termed the "solar composition problem". While the neutrino fluxes were not known to the precision needed to distinguish between the competing models, it had been shown that the solar core metallicity could be extracted directly from future CN solar neutrino measurements, if the precision of associated nuclear cross sections were improved (Haxton and Serenelli, 2008). Reflecting the impact of new experimental work and improved theory, the SF II study of Adelberger *et al.* (2011) significantly revised the SF I cross section for the driving reaction of the CNO-I cycle, and updated key pp-chain cross sections.

The present SF III review began with a four-day workshop in July, 2022, “Solar Fusion Cross Sections III.” The meeting was hosted by Physics Frontier Center N3AS (Network for Neutrinos, Nuclear Astrophysics, and Symmetries)¹, a consortium of institutions involved in multi-messenger astrophysics, whose central hub is the University of California, Berkeley. Sessions were held at the David Brower Center, Berkeley, and the Physics Department, UC Berkeley, with the assistance of N3AS staff. The workshop was organized by N3AS, by Solar Fusion’s long-term sponsor, the INT, and by the European ChETEC-INFRA Starting Community for Nuclear Astrophysics².

The 47 workshop participants represented most of the leading experimental and theoretical research groups that are active in the field. Workshop participants were organized into nine working groups, who were charged with reviewing and evaluating work completed since SF II, to produce updated recommendations. The structure of this review largely follows that of the working groups. The timing of SF III was driven in part by planned large-scale asteroseismic surveys of hydrogen-burning stars such as ESA’s PLATO³ mission (Rauer *et al.*, 2014), and the expectation that associated data analysis will require large libraries of stellar models in which masses, ages, metallicities, and other parameters are varied. SF III will provide the best current nuclear astrophysics input for such modeling. This review summarizes that input, based on the recommendations of the nine SF III working groups.

B. Terminology used

The terminology used here follows recent practice, as summarized for example in the NACRE compilation (Angulo *et al.*, 1999). It is briefly described to aid the reader.

Nuclear reactions are designated as

$$A(b, c)D, \quad (1)$$

where A is the target nucleus, b is the bombarding particle, and c and D are the reaction products, with D distinguished as the heavier product. Frequently in nuclear astrophysics, the reaction cross section σ is re-expressed in terms of the S-factor S_{ij} , with S related to σ as described below, and with i, j the mass numbers of entrance channel nuclei b and A .

The relation between the nuclear cross section $\sigma(E)$ and astrophysical S-factor is

$$S(E) = \sigma(E) E \exp\left(2\pi \frac{Z_A Z_b \alpha}{\beta}\right), \quad (2)$$

with Z_A and Z_b the charge numbers of the target and projectile, $\alpha \sim 1/137$ the fine structure constant, $\beta \equiv v/c = \sqrt{2E/\mu c^2}$ the relative velocity of the interacting particles in units of c , μ the reduced mass, and E the center-of-mass energy. (In this review we use energy units of MeV unless otherwise stated.) This removes the sharp energy dependence from the cross section associated with s-wave Coulomb scattering off a point nucleus, thereby accounting for the leading penetration effects associated with the Coulomb barrier. The quantity $Z_A Z_b \alpha / \beta$ is known as the Sommerfeld parameter. Consequently, unless there are resonances, $S(E)$ varies gently, and at low energies can generally be expressed as a low-order polynomial in E .

The S-factor in Eq. (2) is defined in terms of the nuclear (or “bare”) cross section, while laboratory experiments are performed with targets consisting of neutral atoms. As electrons shield the nucleus, lowering the Coulomb barrier, their presence enhances the laboratory cross section. The S-factors derived from laboratory data must be corrected for this effect. As the energy is lowered, the screening correction becomes increasingly important. Thus particular care must be taken when extrapolating laboratory cross sections to lower energies, as Fig. 4 of SF II illustrates. Unless otherwise noted, the S-factors discussed here are those for bare nuclei.

The rapid rise in the nuclear cross section with energy, reflecting the higher probability of penetrating the Coulomb barrier, competes with a rapid decrease in the probability of finding interacting particles with the requisite center-of-mass energy E in the high-energy tail of the Maxwell-Boltzmann distribution. Consequently, the thermonuclear reaction rate $\langle\sigma v\rangle$ for a given stellar temperature T takes the form

$$\langle\sigma v\rangle = \sqrt{\frac{8}{\mu\pi}} (k_B T)^{-\frac{3}{2}} \int_0^\infty S(E) \exp\left[-\frac{b}{\sqrt{E}} - \frac{E}{k_B T}\right] dE, \quad (3)$$

where $b = \pi Z_A Z_b \alpha \sqrt{2\mu c^2}$ is proportional to the Sommerfeld parameter of Eq. (2) and k_B is Boltzmann’s constant. The maximum of the integrand defines the Gamow peak – the most probable energy for interactions – while the $\sim 1\sigma$ range around the peak is frequently termed the Gamow window.

For the solar reactions of interest, the energy at the Gamow peak is well below the height of the Coulomb potential. The slowly varying $S(E)$ can be expanded as a power series around $E = 0$, with the leading term being $S(0)$, and with corrections given by derivatives taken at $E = 0$,

$$S(E) \simeq S(0) + S'(0) E + \frac{S''(0)}{2} E^2 + \dots \quad (4)$$

Most of the results presented in this review will be given in terms of $S(0)$ and its derivatives.

For most of the nuclear reactions studied here, the notation S_{ij} is adopted, where i and j are the mass numbers

¹ <https://n3as.berkeley.edu/>

² <https://www.chetec-infra.eu/>

³ <https://platomission.com/>

of the projectile and target nuclei, respectively. A summary of SF III reactions and recommended astrophysical S-factors are presented in Table I.

Finally, we provide a short list of abbreviations that are frequently used throughout several sections:

ANC: asymptotic normalization coefficient.

BBN: Big Bang nucleosynthesis.

CME: common-mode error.

EFT: effective field theory, and variants: chiral (χ EFT), pion-less (π EFT), hybrid (EFT*), etc.

LO, NLO, NNLO: leading order, next to (next to) leading order, etc.

SSM: Standard Solar Model.

C. Scope and structure of the review

In the Sun, the primary mechanism for the conversion of four protons into ^4He is the pp chain of Fig. 1, during which two charge-changing weak interactions take place, each converting a proton into a neutron with the emission of a neutrino. The pp chain consists of three main branches, ppI, ppII, and ppIII. These branches are distinguished by their dependence on the solar core temperature and by the neutrinos they produce, with the pp, ^7Be , and ^8B neutrinos serving as “tags” for the ppI, ppII, and ppIII chains, respectively. The SSM predicts that the fluxes of pp, ^7Be , and ^8B neutrinos have an approximate *relative* temperature scaling of $\sim 1:T_c^{10}:T_c^{22}$, where T_c is the solar core temperature (Haxton, 1995). By the mid-1990s, solar neutrino experiments established that the measured fluxes differed significantly from this expected pattern. This contributed to growing expectations that new neutrino physics might be the solution of the solar neutrino problem. It also was an important motivation for SF I, as errors in the nuclear physics of the pp chain could distort the relationship between the various fluxes.

The pp chain accounts for 99% of solar hydrogen burning and are the main focus of this review. The remaining 1% is generated through the CNO-I cycle, see Fig. 2). Motivated in part by Borexino’s recent success in measuring the flux of solar CN neutrinos, this review also considers in more detail than past Solar Fusion reviews the reactions driving hydrogen burning in stars more massive than our Sun.

The review is structured as follows. Section II begins with a summary of past solar neutrino experiments, highlighting some of the open questions that remain. It discusses the second-generation experiments Super-Kamiokande, the Sudbury Neutrino Observatory, and Borexino that contributed in the resolution of the solar neutrino problem. It also describes Hyper-Kamiokande, SNO+, JUNO, and DUNE, new detectors in various stages of construction.

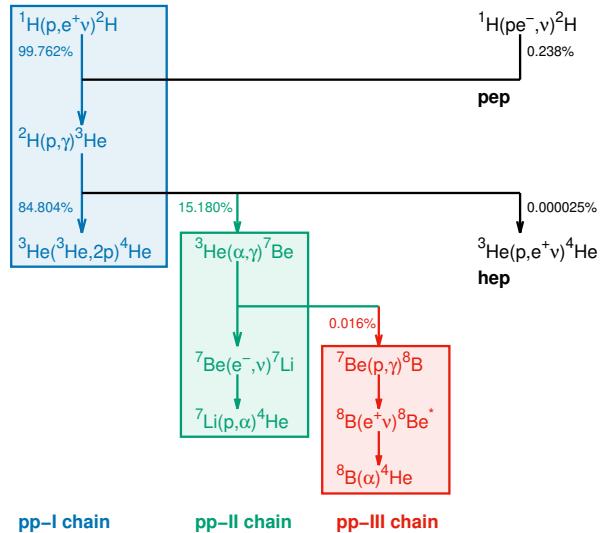


FIG. 1 Nuclear reactions of the proton-proton (pp) chain. The percentage branchings are applicable for the Magg *et al.* (2022) solar composition.

The cross section of the driving reaction of the pp chain, $^1\text{H}(p, e^+\nu)^2\text{H}$, is too small to be measured and thus must be taken from theory. This important reaction is reviewed in section III. With the exception of the minor hep branch, all other reactions and decays have been studied in laboratories, with measurements then combined with theory to predict rates under solar conditions. The $^2\text{H}(p, \gamma)^3\text{He}$, $^3\text{He}(^3\text{He}, 2p)^4\text{He}$, and $^3\text{He}(\alpha, \gamma)^7\text{B}$ reactions are reviewed in Secs. IV, V, and VI, respectively. The hep reaction is reviewed in Section VII, and the electron-capture reactions on pp and ^7Be are discussed in Section VIII. The $^7\text{Be}(p, \gamma)^8\text{B}$ reaction responsible for the high energy neutrinos measured in Super-Kamiokande is treated in Section IX.

While a modest contributor to solar energy production, the CNO-I cycle is a potentially important probe of core metallicity as its hydrogen burning is catalyzed by the Sun’s primordial C and N (Section XI).

A current focus of nuclear astrophysics laboratories is the higher-temperature pathways for hydrogen burning that operate in massive stars. Many of the contributing reactions have been re-measured. That progress is reviewed in Section XI.

Charged-particle nuclear reactions occurring at energies below the Coulomb barrier are affected by electron screening, the shielding of the nuclear charge by electrons. Because of their very different atomic environments, laboratory reactions and those occurring in the solar plasma are affected in distinct ways. Section XII describes the current status of efforts to account for these differences, when solar cross sections are extracted from laboratory measurements.

The energy transport inside the radiative zone is af-

TABLE I List of nuclear reactions reviewed in SF III. Denoting the astrophysical S-factor by S_{ij} , its value at zero-energy $S(0)$ is given along with, where applicable, derivatives parameterized in Eq. (4). See the corresponding section for uncertainties, higher precision values, and detailed discussion.

Reaction	S_{ij}	$S(0)$ (MeV b)	$S'(0)$ (b)	$S''(0)$ (MeV $^{-1}$ b)	Section
$^1\text{H}(p, e^+\nu)^2\text{H}$	S_{11}	4.09×10^{-25}	4.5×10^{-24}	9.9×10^{-23}	III
$^2\text{H}(p, \gamma)^3\text{He}$	S_{12}	2.03×10^{-7}	see text		IV
$^3\text{He}(^3\text{He}, 2p)^4\text{He}$	S_{33}	5.21	-4.90	22.42	V
$^3\text{He}(\alpha, \gamma)^7\text{Be}$	S_{34}	5.61×10^{-4}	-3.03×10^{-4}	-	VI
$^3\text{He}(p, e^+\nu)^4\text{He}$	S_{hep}	8.6×10^{-23}	-	-	VII
$^7\text{Be}(p, \gamma)^8\text{B}$	S_{17}	2.05×10^{-5}	-	-	IX
$^{14}\text{N}(p, \gamma)^{15}\text{O}$	S_{144}	1.68×10^{-3}	-	-	X.F
$^{12}\text{C}(p, \gamma)^{13}\text{N}$	S_{112}	1.44×10^{-3}	2.71×10^{-3}	3.74×10^{-2}	XI.A.1
$^{13}\text{C}(p, \gamma)^{14}\text{N}$	S_{113}	6.1×10^{-3}	1.04×10^{-2}	9.20×10^{-2}	XI.A.2
$^{15}\text{N}(p, \gamma)^{16}\text{O}$	S_{115}^γ	4.0×10^{-2}	1.07×10^{-1}	1.84	XI.B.1
$^{15}\text{N}(p, \alpha)^{12}\text{C}$	S_{115}^α	73	3.37×10^2	1.32×10^4	XI.A.3
$^{16}\text{O}(p, \gamma)^{17}\text{F}$	S_{116}	1.09×10^{-2}	-4.9×10^{-2}	3.11×10^{-1}	XI.B.2
$^{17}\text{O}(p, \gamma)^{18}\text{F}$	S_{117}	4.7×10^{-3}	-	-	XI.B.3
$^{18}\text{O}(p, \gamma)^{19}\text{F}$	S_{118}	2.30×10^{-2}	-	-	XI.C.1
$^{20}\text{Ne}(p, \gamma)^{21}\text{Na}$	S_{120}	6.78	-	-	XI.D.1
$^{21}\text{Ne}(p, \gamma)^{22}\text{Na}$	S_{121}	$\approx 2.0 \times 10^{-2}$	-	-	XI.D.2
$^{22}\text{Ne}(p, \gamma)^{23}\text{Na}$	S_{122}	0.415	-	-	XI.D.3
$^{23}\text{Na}(p, \gamma)^{24}\text{Mg}$	S_{123}	1.80×10^{-2}	0	0	XI.D.4

fected by the metals⁴ that were incorporated into the Sun when it first formed. While metals comprise less than 2% of the Sun by mass, they play an outsized role in determining the opacity. Very few of the needed radiative opacities have been measured, and the conditions under which measurements are made typically are not identical to those in the Sun. Section XIII describes the current state of the art, and discusses connections between opacities and solar composition that have complicated efforts to resolve the solar metallicity problem.

Much of the progress made in constraining the reactions of the pp chain and CNO-I cycles has been possible thanks to new laboratories and the facilities they host. The underground LUNA laboratory is a prominent example (Broggini *et al.*, 2018). Section XIV describes the capabilities of selected laboratories involved in the measurements reviewed here.

II. Solar neutrino observations

Solar neutrinos offer a unique tool box for probing both the fundamental properties of these elusive particles, and their interactions with matter, as well as understanding their source: the fusion reactions that power our Sun. The original motivation for observations of solar neutrinos was precisely the hope to probe solar fusion. The first

successful experiment beginning in 1967 at the Homestake mine (Cleveland *et al.*, 1998) offered the surprising result of a neutrino flux suppressed to approximately one third of expectation. The intervening years have seen a tour de force of experimental efforts, leading to the resolution of the Solar Neutrino Problem and the confirmation of neutrino oscillations (Ahmad *et al.*, 2002a, 2001; Ahmed *et al.*, 2004; Eguchi *et al.*, 2003; Fukuda *et al.*, 1998). These high-precision flux and spectral measurements mapped out the details of solar fusion, probed the structure of the Sun, and demonstrated (together with atmospheric neutrino measurements) that neutrinos have masses and mix. Table II lists the pp-chain and CNO cycle neutrino sources, their endpoint (β decay) or line (electron capture) energies E_ν , and predicted SSM fluxes.

A. Open questions

The solar neutrino spectrum has been measured for energies above 233 keV, with some fluxes determined to a precision equivalent to or better than theoretical predictions. However, improved measurements can offer further insight on several aspects. For instance, more precise determinations of the fluxes of neutrinos from the CNO cycle (Fig. 2) could determine the contemporary abundances of N and O in the solar core (Castro *et al.*, 2007; Kunitomo *et al.*, 2022; Serenelli *et al.*, 2011). While these abundances differ significantly from the primordial values that are input into the SSM –the primordial abundances are altered by out-of-equilibrium CNO-cycle burning that occurs in the early Sun (Roxburgh, 1985)– this burning

⁴ In this work, we follow the astrophysical terminology where "metals" refer to all chemical elements except for the two lightest ones, hydrogen and helium.

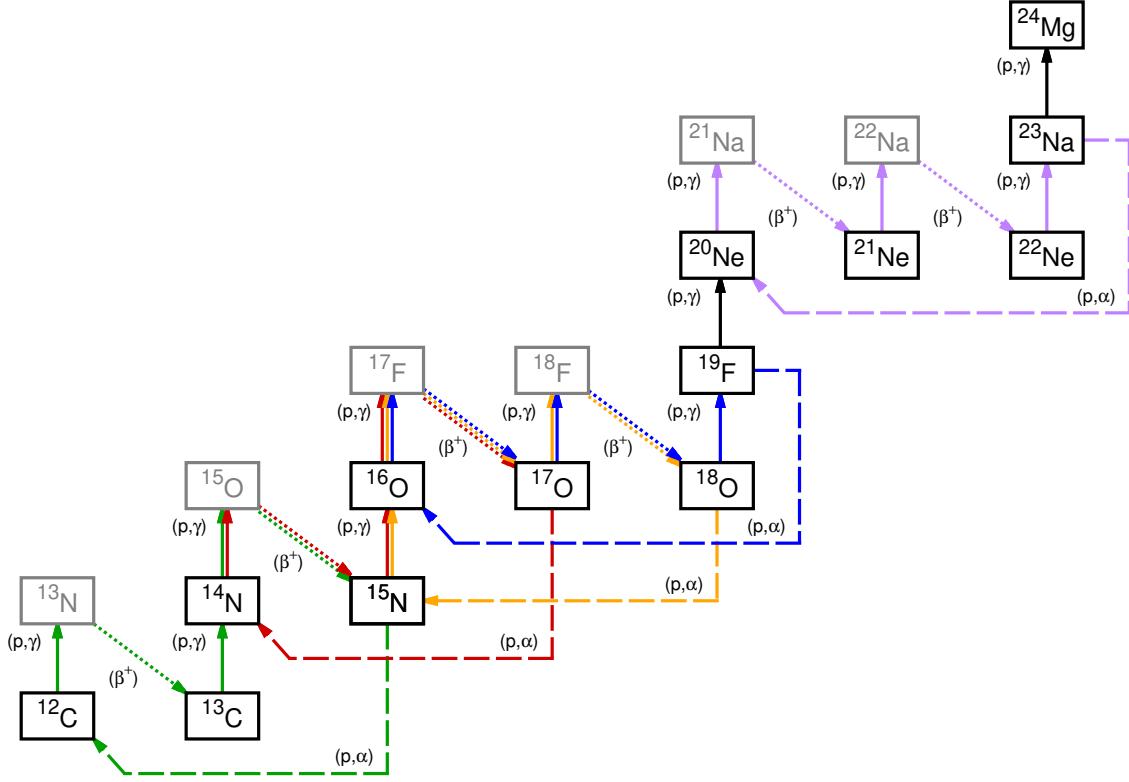


FIG. 2 Nuclear reactions of the CNO-I (green), CNO-II (red), CNO-III (orange), CNO-IV (blue), and Ne-Na (purple) cycles. Other reactions reported in this work are also present (in black). For visual aid, (p, γ) reactions are styled using full arrows, while (p, α) reactions use dashed-line arrows. Greyed isotopes are short-lived and quickly (β^+) -decay (dotted arrows), resulting in $e^+ + \nu_e$ emissions.

TABLE II Solar neutrino sources, energies, and SSM flux predictions. All β decay sources produce continuous spectra, while the pep and ${}^7\text{Be}$ electron-capture sources produce line spectra. The fluxes are taken from the SSM calculations of Herrera and Serenelli (2023), computed with nuclear reaction rates shown in Table I, for the compositions of GS98 (Grevesse and Sauval, 1998) (high-Z), AAG21 (Asplund *et al.*, 2021) (low-Z), and MB22p (Magg *et al.*, 2022) (high-Z) with associated uncertainties indicated.

Source	E_ν (MeV)	GS98	AAGS21	MB22p	units $(\text{cm}^{-2}\text{s}^{-1})$
pp	≤ 0.420	5.96 (0.6%)	6.00 (0.6%)	5.95 (0.6%)	10^{10}
pep	1.442	1.43 (1.1%)	1.45 (1.1%)	1.42 (1.3%)	10^8
hep	≤ 18.77	7.95 (31%)	8.16 (31%)	7.93 (30%)	10^3
${}^7\text{Be}$	0.862 ^(a) 0.384 ^(b)	4.85 (7.4%)	4.52 (7.3%)	4.88 (8.1%)	10^9
${}^8\text{B}$	$\lesssim 15$	5.03 (13%)	4.31 (13%)	5.07 (15%)	10^6
${}^{13}\text{N}$	≤ 1.198	2.80 (16%)	2.22 (13%)	3.10 (15%)	10^8
${}^{15}\text{O}$	≤ 1.732	2.07 (18%)	1.58 (16%)	2.30 (18%)	10^8
${}^{17}\text{F}$	≤ 1.738	5.35 (20%)	3.40 (16%)	4.70 (17%)	10^6

Notes: (a) 90% and (b) 10% of the ${}^7\text{Be}$ neutrino flux, respectively.

redistributes the C, N, and O, but does not alter the net

number of such nuclei. Consequently, CNO neutrino flux measurements might allow one to experimentally address the solar composition problem discussed below.

Solar neutrino flux measurements can also constrain physical processes in stars such as chemical mixing. The fluxes are sensitive to solar/stellar model inputs, including the radiative opacities (see Sect. XII). Measurements of the neutrino fluxes from ${}^8\text{B}$ and ${}^7\text{Be}$ can offer a handle on the temperature of the solar core, and other environmental factors (Bahcall and Ulmer, 1996).

A measurement of the pp flux with percent-level precision would provide a test of the luminosity constraint, further probing solar power generation mechanisms (Bahcall, 2002; Vescovi *et al.*, 2021). The one branch of the pp chain that yet remains undetected, the hep flux, has both the highest energy but also the lowest flux of neutrinos from the Sun. The spectrum extends beyond that of the ${}^8\text{B}$ neutrinos, offering a small window for unequivocal observation. Limits have been placed on this flux (Aharmim *et al.*, 2020), but a definitive observation will likely require a large, next-generation detector such as Hyper-Kamiokande (HyperK) (Hyper-Kamiokande Proto-Collaboration *et al.*, 2018) or the Deep Underground Neutrino Experiment

(DUNE) (Capozzi *et al.*, 2019).

Finally, the observation of monochromatic neutrinos produced by electron capture reactions on ^{13}N , ^{15}O and ^{17}F (which we refer to as ecCNO neutrinos) (Bahcall, 1990; Stonehill *et al.*, 2004; Villante, 2015) could be rewarding in terms of physical implications. Indeed, they can be used as probes of the metallicity of the solar core. Moreover, they provide a measure of the electron neutrino survival probability at specific neutrino energies $E_\nu \sim 2.5$ MeV within the broad transition region between vacuum and matter-enhanced oscillations, where currently we have no constraints. This would constitute a new test of the large-mixing-angle (LMA) MSW flavor oscillation paradigm. The detection of this subdominant component of the CNO-cycle is extremely difficult but it could be within reach of future very large ultra-pure liquid scintillator detectors (Villante, 2015).

Solar neutrinos provide an opportunity to understand the interaction of neutrinos with matter. The effect of MSW oscillations has a significant impact on the observed solar neutrino spectrum: with vacuum oscillation dominating at low energy, below approximately 1 MeV, where the survival probability is roughly one half, while matter effects further suppress the flux to a survival probability of roughly one third above approximately 5 MeV. The transition region between these two regimes offers an extremely sensitive probe of the details of the interactions of neutrinos with matter, including the potential to search for new physics such as sterile neutrinos, or non-standard interactions, by looking for distortions to the expected spectral shape (Bhupal Dev *et al.*, 2019; Borexino Collaboration *et al.*, 2020a; Coloma *et al.*, 2023).

The same theory of matter effects predicts a small regeneration of electron neutrinos during the night time as they propagate through the bulk of the earth. This so-called “day/night effect” has been sought after by both Super-Kamiokande (SuperK) and the Sudbury Neutrino Observatory (SNO) (Ahmad *et al.*, 2002b; Renshaw *et al.*, 2014), but a significant observation is still limited by statistics. Future data from HyperKamiokande or DUNE may be needed to confirm our understanding of this prediction of the MSW effect.

The slight tension between measurements of the mass splitting parameters, Δm_{12}^2 in solar neutrino experiments and terrestrial data sensitive to the same parameter, from the KamLAND reactor experiment has now disappeared with the addition of new data on day-night effects and spectral shapes (Capozzi *et al.*, 2018; Esteban *et al.*, 2020). Further tightening of the oscillation parameters can be expected from JUNO, which will obtain highly precise data on both the solar mass splitting and the mixing angle (An *et al.*, 2016).

1. The solar composition problem

Solar photospheric abundances, determined with spectroscopic techniques, are a fundamental input for the construction of SSMs, but also they are used as input to nearly all models in astrophysics, including stellar evolution, protoplanetary disks, and galactic chemical evolution. Over the last two decades, the development of three dimensional radiation hydrodynamic models of the solar atmosphere (Asplund *et al.*, 2000; Ludwig *et al.*, 2009) and of techniques to study line formation under non-local thermodynamic conditions (Asplund, 2005), together with improved descriptions of atomic properties (e.g. transition strengths Bergemann *et al.* 2021; Magg *et al.* 2022), have led to a significant revision of solar abundances. Initial results based on these more sophisticated methods favored a markedly lower solar metallicity (Asplund *et al.*, 2009), particularly in the CNO elements, than obtained in the 90s from older techniques (Grevesse and Noels, 1993; Grevesse and Sauval, 1998).

Recently, two groups have revisited the solar photospheric abundances using modern methods (Asplund *et al.*, 2021; Magg *et al.*, 2022). While Asplund *et al.* (2021) obtain results consistent with their previous findings, Magg *et al.* (2022) find an O abundance intermediate between that of Asplund’s group and those from the 90s, in agreement with another 3D-based determination by Caffau *et al.* (2011). Interestingly, Magg *et al.* (2022) find higher C and N abundances and, indirectly, a higher Ne abundance due to the larger Ne to O ratio measured in the solar corona (Young, 2018). This leads to a combined metal-to-hydrogen ratio that is by chance comparable to those from Grevesse and Noels (1993) and Grevesse and Sauval (1998), albeit with a different mixture of elements.

Considering that uncertainties in element abundances are difficult to quantify, it has become customary to consider two canonical sets of abundances, which we refer to as high metallicity (HZ) and low metallicity (LZ) solar admixtures. See, for example, Vinyoles *et al.* (2017) for SSM reference values. In this context, the new solar abundance determinations by Asplund *et al.* (2021) fall into the LZ category, while those from Magg *et al.* (2022) are HZ. Solar models employing the LZ abundances fail to reproduce most helioseismic probes of solar properties. This disagreement constitutes the so-called solar composition problem (Bahcall *et al.*, 2005a; Basu and Antia, 2004; Delahaye and Pinsonneault, 2006) that has defied a complete solution. All proposed modifications to physical processes in SSMs offer, at best, only partial improvements in some helioseismic probes, see e.g. (Basu and Antia, 2008; Castro *et al.*, 2007; Guzik and Mussack, 2010; Guzik *et al.*, 2005; Serenelli *et al.*, 2011). The same conclusions are obtained with SSMs computed with the newest LZ (Asplund *et al.*, 2021) and HZ (Magg *et al.*, 2022) abundances as discussed also there. An al-

ternative possibility is to consider modifications to the physical inputs of SSMs at the level of the constitutive physics, radiative opacities in particular. This is possible because most helioseismic probes depend actually not directly on the solar composition, but on the radiative opacity profile in the solar interior, i.e. on the combination of solar composition and atomic opacities (see § XIII). The same can be said for solar neutrinos from the pp chain. Early work (Bahcall *et al.*, 2005b; Montalbán *et al.*, 2004) already suggested that a localized increase in opacities could solve or, at least, alleviate the disagreement of low-Z solar models with helioseismology, and Christensen-Dalsgaard *et al.* (2009) and Villante (2010) showed that a tilted increase in radiative opacities, with a few percent increase in the solar core and a larger (15–20%) increase at the base of the convective envelope could lead to LZ SSMs that would satisfy helioseismic probes equally well as HZ SSMs.

The degeneracy between solar composition and opacities can be broken using CNO solar neutrinos, e.g., following the methodology developed in Haxton and Serenelli (2008). Such a study was recently carried out by the Borexino collaboration (Appel *et al.*, 2022; Basilico *et al.*, 2023; Borexino Collaboration *et al.*, 2020b). The Borexino measurement of the CNO-I cycle neutrino flux (the ^{13}N and ^{15}O fluxes) were used to determine the C+N core abundance. Results show a $\sim 2\sigma$ tension with LZ metallicity determinations, while being in better agreement with HZ mixtures. While the error budget is presently dominated by the uncertainty of the Borexino CNO neutrino measurement, a significant contributor to the error ($\sim 10\%$) is nuclear, due to uncertainties in S_{114} , S_{34} , and S_{17} . The interpretation of future improved CNO neutrino measurements will be impacted, unless these nuclear physics uncertainties are reduced.

Very recently, Gonzalez-Garcia *et al.* (2024) presented a new global determination of all solar neutrino fluxes using all available experimental data, including the latest phases of Borexino. Results from this global analysis are in line with those from Borexino, although the added ^{13}N and ^{15}O fluxes are about 10% lower than the Borexino result alone. A comparison of the solar fluxes with SSM calculations (Herrera and Serenelli, 2023) shows that HZ SSMs are in better agreement with solar neutrino fluxes than the LZ SSMs, pointing toward a C+N solar core abundance consistent with HZ abundances. The discrimination that solar neutrino fluxes can offer between solar compositions is at most, however, of the order of 2σ . (See Table 2 in Gonzalez-Garcia *et al.* (2024)).

Helioseismology also offers the potential to determine the total solar metallicity, i.e. without disentangling individual element abundances, rather independently from opacities. This relies on using the so-called adiabatic index Γ_1 , which deviates from $5/3$ in regions of partial ionization. The underlying technique has been used widely to determine the helium abundance in the solar envelope

and it has also been extended to determine solar metallicity. Attempting the latter is difficult because the imprint of partial ionization of metals is quite subtle. Previous work along this line (Antia and Basu, 2006) found an overall metallicity consistent with HZ abundances, but new work (Buldgen *et al.*, 2024) claims to favor LZ values. While this method depends very weakly on radiative opacities, the abundance determination is degenerate with the equation of state. Independent confirmation of these results would be desirable. In short, the controversy related to the solar composition is far from being resolved, currently with different indicators showing contradictory results.

2. The Gallium anomaly

With the aid of very intense radioactive sources of ^{51}Cr and ^{37}Ar , tests have been made of the rate of production of ^{71}Ge by neutrino interactions on ^{71}Ga , the basis of radiochemical measurements of the low-energy solar neutrino flux. Initial experiments showed lower rates than expected, with interesting but inconclusive statistical precision. Very recently the BEST experiment (Barnov *et al.*, 2018, 2022a,b) has confirmed this anomaly at more than 4σ . Many possible explanations have been explored, but for each there are contradictions (Akhmedov and Smirnov, 2022; Brdar *et al.*, 2023; Giunti *et al.*, 2023; Krueger and Schwetz, 2023). A sterile neutrino explanation is disfavored (Giunti *et al.*, 2022; Goldhagen *et al.*, 2022), particularly because of conflict with solar neutrino limits (see Fig. 3), as well as cosmological bounds (Hagstotz *et al.*, 2021).

The sterile neutrino contribution to the solar flux is fundamentally limited by the luminosity constraint, that the Sun's total energy output is the result of nuclear reactions that produce neutrinos, whether active or sterile. If precise neutral-current data were available across the solar spectrum, a completely model-independent limit on a possible sterile component could be determined. The present neutral-current data at low energies (from Borexino) are imprecise but can be supplemented by much more precise charged-current data together with 3-flavor oscillation physics. Still better limits, at the cost of some model dependence, can be obtained with fits of experimental data to solar models, as shown in Fig. 3 above, as well as Fig. 1 of Goldhagen *et al.* (2022).

The precisely known rate of electron capture on ^{71}Ge to ^{71}Ga , for which the half-life is $11.43(3)$ d (Hampel and Remsberg, 1985), places an important constraint on the neutrino absorption rate (Elliott *et al.*, 2023; Giunti *et al.*, 2023). The half-life has now been remeasured by two groups, Norman *et al.* (2024) and Collar and Yoon (2023), yielding values of $11.468(8)$ d and $11.46(4)$ d, consistent with the previous determination while increasing the precision. The allowed matrix element that governs

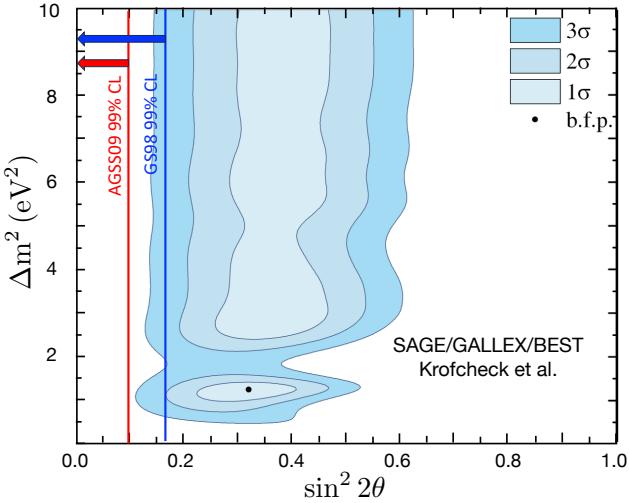


FIG. 3 The contours are the sterile neutrino interpretation of the Ga anomaly, from (Elliott *et al.*, 2023). Limits at the 99% C.L. on the mixing angle from solar neutrinos, from (Goldhagen *et al.*, 2022), are indicated at left for two solar models, GS98 and AGSS09, and essentially exclude the indicated sterile-neutrino space.

neutrino capture to the ground state of ^{71}Ge can be extracted to a precision of $\sim 1\%$ from the electron capture rate. The ground-state transition by itself generates a significant anomaly: inclusion of the excited-state contributions doubles the effect. In Appendix A we describe some of the details of both the 1997 extraction of the allowed (Gamow Teller) strength by Bahcall (1997) and the recent update of Elliott *et al.* (2024, 2023).

Elliott *et al.* (2024, 2023) find the ground-state cross section for absorbing ^{51}Cr neutrinos is 2.5% lower than Bahcall's value. This work includes the contributions of weak magnetism and radiative corrections, which are shown to be sub-1% effects. An improved extraction of the $\sim 6\%$ excited-state contribution to neutrino absorption was performed using data from forward-angle (p,n) scattering. This led to a slight increase in that contribution relative to Bahcall (1997), thereby reducing the 2.5% difference above by about half, for the total ^{51}Cr and ^{37}Ar capture cross sections.

The procedures followed by Elliott *et al.* (2024, 2023) could be extended to the solar neutrino ^{71}Ga capture cross section, including the contribution from the high-energy ^8B neutrinos. That has not been done, but as noted in Appendix A, could help resolve small discrepancies noted there.

Results from the SAGE and GALLEX/GNO experiments remain part of the solar neutrino database used in various global fits to neutrino parameters. A resolution of the anomaly that emerged from BEST and the four gallium detector calibration experiments is important as it would increase confidence in these data.

3. The Boron-8 neutrino spectrum

The neutrino spectrum from ^8B , extending to approximately 17 MeV, has been the most accessible part of the solar neutrino spectrum (Abe *et al.*, 2016; Aharmim *et al.*, 2020), playing an important role in disclosing the physics of neutrino oscillation and testing the SSM. Knowing it with precision, particularly at the high energy end, has renewed interest in the context of observing the hep neutrinos (Aharmim *et al.*, 2020, 2006), which could be accessible with future observatories (Askins *et al.*, 2022).

In SF II it was decided to recommend the spectrum calculated by Winter et al. (Winter, 2007; Winter *et al.*, 2003, 2006) based on their measurement of the alpha spectrum from ^8B . This spectrum showed excellent agreement with an independent experiment (Bhattacharya *et al.*, 2006). Both radiative and recoil-order corrections were included by Winter et al. The radiative corrections are relatively small, due to a cancellation between the real and virtual contributions (Batkın and Sundaresan, 1995). The recoil-order corrections are dominated by the weak magnetism part, which has been deduced from measurements of the analog electromagnetic decays (de Braeckeleer *et al.*, 1995).

Since then three additional measurements (Kirsebom *et al.*, 2011; Longfellow *et al.*, 2023; Roger *et al.*, 2012) of the ^8B β -decay alpha spectrum have been performed. All three find that the peak of the alpha spectrum appears about 20 keV lower than determined by Winter et al. Overall, these new measurements yield differences in the ^8B neutrino spectrum of $\lesssim 5\%$ below $E_\nu = 15$ MeV, the energy above which the ^8B contribution becomes small compared to the hep component. Figure 4 shows the ^8B spectral uncertainties on the scale of the hep spectrum. With resolution effects considered, it will be difficult to extract the hep flux from solar neutrino measurements without further reduction in ^8B uncertainties.

In producing a recommendation for the ^8B neutrino spectrum we considered the following:

1. Apparently the uncertainty estimations for the neutrino spectrum in Roger *et al.* (2012) are not quite correct. Longfellow *et al.* (2023) show that the uncertainty due only to the weak magnetism part, for which all authors follow similar prescriptions, is larger in the 0–12 MeV range than the overall uncertainty estimated by Roger et al. We estimated the uncertainties due to the weak magnetism term only and found agreement with those of Longfellow et al.
2. All three of the recent efforts (Kirsebom *et al.*, 2011; Longfellow *et al.*, 2023; Roger *et al.*, 2012) used detectors (Si strip) more complicated than those used by Winter et al. (a Si surface barrier

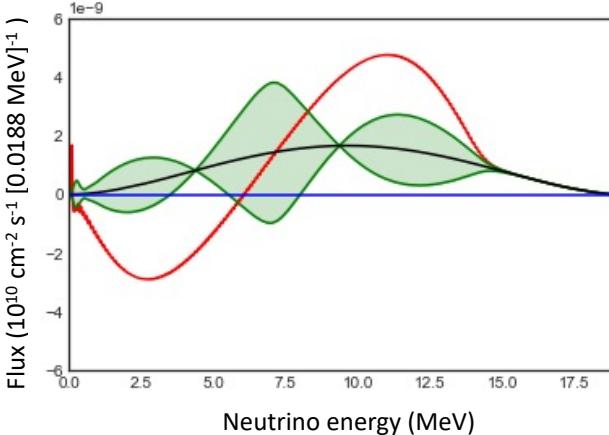


FIG. 4 The hep spectrum is in black, and the uncertainties in ${}^8\text{B}$ from Longfellow *et al.* (Longfellow *et al.*, 2023) span the region between the green lines. The red line indicates the difference between Longfellow *et al.* and Winter *et al.* (Winter *et al.*, 2006) deduced ${}^8\text{B}$ spectra. The hep spectrum is used here as a convenient metric for the size of the uncertainties in the ${}^8\text{B}$ spectrum.

detector). The strip detectors allow for better position resolution and reduced beta summing. However, for detection from sources external to the detectors (as opposed to detection from sources implanted deep into a detector), the complicated pattern of dead layers needs to be measured and taken into consideration for accurate calibrations. Effects of partial charge collection in-between strips are also significant. The effects are different for a calibration source than from the emission from ${}^8\text{B}$ trapped as an ion.

Based on these considerations, we recommend the spectrum of Longfellow *et al.* (2023), but suggest that conclusions sensitive to the choice of ${}^8\text{B}$ neutrino spectrum, such as evidence for hep neutrinos, should be evaluated using both the Winter *et al.* and the Longfellow *et al.* spectra.

Given the importance of the weak magnetism contributions, it would be good to have a new experiment with reduced uncertainties that could be compared directly to the results of de Braeckeleer *et al.* (1995). Recoil-order effects in the $A = 8$ system were most recently measured by Sumikama *et al.* (2011), making use of the alignment- β correlation. Using combinations of their measurements and $\beta - \alpha$ correlation data from McKeown *et al.* (1980), they were able to extract some of the recoil-order matrix elements. Calculations of these matrix elements in the symmetry-adapted no-core shell model were recently performed by Sargsyan *et al.* (2022). The comparison to the experimentally derived matrix elements of McKeown *et al.* (1980), de Braeckeleer *et al.* (1995), and Sumikama *et al.* (2011) is not straightforward because the experiments quote averages over the whole beta spectrum as

opposed to fits for each of the four levels used in the calculation. A more detailed comparison of theory to experiment that takes into account the averaging would be helpful. An additional point of comparison that has been used previously is the beta spectrum from ${}^8\text{B}$ (Bahcall *et al.*, 1996). An experiment with improved statistics and systematics that extends over the full range of beta energies would be very useful. The existing data from Napolitano *et al.* (1987) is in good agreement with the alpha spectrum from Winter *et al.* But given the apparent 20 keV difference that the newer alpha-spectrum experiments described above report, it would be good to have a modern beta spectrum measurement specifically designed to address this difference.

B. Experimental program on solar neutrinos

A broad range of technology can be used to interrogate solar neutrinos. Radiochemical experiments utilizing chlorine and gallium played a crucial role in the early days of the solar neutrino problem. Large monolithic detectors have achieved great success in real-time measurements – from the water Cherenkov detectors such as Kamiokande, SNO, and SuperK, which could use the unique topology of Cherenkov light to point back to the neutrinos’ origin, to liquid scintillator detectors such as KamLAND and Borexino, whose high light yield and low threshold allow for precision spectroscopy.

1. Super-Kamiokande and Hyper-Kamiokande

The first real-time detection of solar neutrinos was achieved by the Kamiokande experiment (Hirata *et al.*, 1989), which detected the neutrinos via elastic scattering of electrons, a process in which the outgoing electron’s direction is highly correlated to that of the incoming neutrino. This allowed Kamiokande to directly point back to the neutrinos’ origin: our Sun. The successor experiment, Super-Kamiokande, is a 50-kton volume of pure water, surrounded by over 10,000 photon detectors, which has achieved an unparalleled program of neutrino observations and other physics over its several decades of operation. This program has included the highest-precision measurement of the elastic scattering signal from solar neutrino interactions, as well as sensitive searches for the day/night distortions of the spectral shape (Abe *et al.*, 2016). Now filled with a gadolinium additive to enhance neutron capture, SuperK’s primary current focus is the search for the Diffuse Supernova Neutrino Background (DSNB) via inverse beta decay (Harada *et al.*, 2023). It also continues its atmospheric neutrino program: SuperK measurements made 25 years ago demonstrated the neutrino oscillations were responsible for the puzzling zenith-angle-dependence of

this flux (Fukuda *et al.*, 1998).

In parallel to ongoing operation of SuperK, an even larger sister project is under construction. At over 250 kton in total mass, and with improved light collection relative to SuperK, Hyper-Kamiokande (HyperK) will impact statistics-limited searches the involve the higher energy ${}^8\text{B}$ solar neutrinos, including day/night effects, the search for hep neutrinos, and the shape of the ${}^8\text{B}$ neutrino spectrum. After 10 years of operations, HyperK will reach a sensitivity to day/night effects $\gtrsim 4$ (8) σ , given the oscillation parameters deduced from reactor (solar) experiments. HyperK will serve as the far detector for JPARC’s long-baseline neutrino program.

2. Sudbury Neutrino Observatory and SNO+

SNO’s unique use of heavy water as a target medium offered the ability to detect solar neutrinos via two additional interactions besides elastic scattering. In the charged current interaction an electron neutrino interacts with the neutron in the deuteron, producing an electron and a proton. This interaction path is sensitive only to electron-flavor neutrinos at the few-MeV energy scale of solar neutrinos, providing a measurement of the pure ν_e flux. SNO also had access to the neutral current interaction, in which a neutrino of any flavor interacts with the deuteron, breaking it apart into its constituent nucleons. Being equally sensitive to all active flavors, the neutral current interaction yields a measurement of the total, flavor-blind neutrino flux. It was this capability that allowed SNO to resolve the solar neutrino problem, demonstrating that most of the ν_e s produced by the Sun were arriving at earth in a different flavor state (Ahmed *et al.*, 2002a, 2001; Ahmed *et al.*, 2004).

SNO ceased data taking in 2006. Analysis of the data set has continued, and produced a number of new results, including constraints on non-standard effects such as Lorentz violation and neutrino decay (Aharmim *et al.*, 2018, 2019a), limits on the fluxes of hep neutrinos and the diffuse supernova neutrino background (DSNB) (Aharmim *et al.*, 2020), and neutron production from cosmogenic muons and from atmospheric neutrino interactions (Aharmim *et al.*, 2019b,c).

After decommissioning, the detector was repurposed for the SNO+ experiment, in which the target material was replaced with a pure organic liquid scintillator: linear alkyl benzene (LAB), loaded with 2.2 g/L of the secondary fluor, PPO. The high light yield of this scintillator, coupled with the location in SNOLAB –at 6-km water equivalent, one of the deepest underground laboratories in the world– and an extremely well-understood detector, allow for a range of high-precision measurement programs. In a preliminary water phase, SNO+ demonstrated detection of neutron captures on hydrogen (Anderson *et al.*, 2020), an impressive technical achievement

in an unloaded water detector; a low background measurement of ${}^8\text{B}$ solar neutrinos (Anderson *et al.*, 2019b); several searches for invisible modes of nucleon decay (Allega *et al.*, 2022; Anderson *et al.*, 2019a); and detection of antineutrinos from reactors over 200 km away (Allega *et al.*, 2023). The low backgrounds achieved and the detector’s demonstrated technical capabilities set the stage for future measurements. In a “partial-fill” stage, in which the upper half of the detector was filled with liquid scintillator while the lower half still contained water, the first demonstration of event-level direction reconstruction was achieved for ${}^8\text{B}$ neutrinos in a scintillator detector (Allega *et al.*, 2024). This was facilitated by a lower loading of PPO at that time, which results in lower scintillation yield and a slower time profile, effectively enhancing the Cherenkov component, which can be leveraged for directional information.

Now fully filled with liquid scintillator, SNO+ is the deepest, largest operating liquid scintillator detector in the world. The future program will include measurements of several solar neutrino fluxes, as well as the ${}^8\text{B}$ neutrino energy spectrum.

3. Borexino

Water Cherenkov experiments such as Kamiokande, SuperK and SNO provided the first real-time measurement of solar neutrinos but their relatively high energy threshold made them sensitive only to a small fraction of the solar neutrino flux. To study in real time the bulk of solar neutrino emission a different detector technology was required. The Borexino experiment, located deep underground at the INFN Laboratori Nazionali del Gran Sasso, used an organic liquid scintillator target made of pseudocumene with 1.5 g/L PPO to detect the elastic scattering of solar neutrinos off electrons. The high light yield of the liquid scintillator made it possible to significantly lower the energy threshold, but given the small signal rate and the lack of a clear signature to separate it from the background (such as the direction indicated by the emission of Cherenkov radiation), the experiment required a long preparatory phase to develop the most advanced techniques to suppress the background, especially the one due to radioactive contamination of the liquid scintillator itself.

The detector’s extreme radiopurity was key to its success, and over its 14 years of data taking, ending in October 2021, Borexino proved itself capable of covering the entire solar neutrino spectrum. At the beginning of its data taking in 2007 the level of ${}^{238}\text{U}$ and ${}^{232}\text{Th}$ contamination in Borexino were lower than 2×10^{-17} and 7×10^{-18} g/g respectively (BOREXINO Collaboration *et al.*, 2008), paving the way for the first measurement of the ${}^7\text{Be}$ sub-MeV solar neutrinos, followed by a low-threshold measurement of the ${}^8\text{B}$ flux and a first indica-

tion of the pep neutrinos. After a purification campaign that further reduced the liquid scintillator contamination, Borexino aimed at improving the accuracy of its results in the measurement of the pp-chain solar neutrinos. The flux of neutrinos produced in the pp fusion process was first reported in [BOREXINO Collaboration et al. \(2014\)](#) and later improved, along with all the other fluxes from the pp chain, apart from the hep neutrinos ([Agostini et al., 2020b, 2019; Borexino Collaboration et al., 2018](#)). The pp flux was determined to an uncertainty of 11%, improving the neutrino-based estimate of solar luminosity, while ^7Be neutrinos were measured with an uncertainty of 2.7%, half that of the SSM prediction. Furthermore, using the SSM to constrain the flux of CNO neutrinos, the pep neutrino signal was established with a significance exceeding 5σ for the first time, and the ^8B flux was measured with a threshold as low as 3 MeV.

After completing the investigation of the pp-chain neutrinos, Borexino reported the first detection of solar neutrinos produced in the CNO-cycle ([Borexino Collaboration et al., 2020b](#)), demonstrating directly that this mode of hydrogen burning operates in stars. Profiting from a larger exposure and a better understanding of the radioactive backgrounds enabled by the unprecedented thermal stability of the detector, Borexino further improved its measurement of the CNO neutrino flux ([Appel et al., 2022; Basilico et al., 2023](#)), where the new CNO result was used in combination with existing solar neutrino results to probe solar composition.

Borexino’s physics program was not limited to solar neutrinos: its outstanding radiopurity made it a excellent detector for geoneutrinos ([Agostini et al., 2020a](#)) and for searches for various rare processes. Borexino data were used to constrain exotic properties of neutrinos, such as their magnetic moments ([Agostini et al., 2017](#)) and non standard interactions with matter ([Borexino Collaboration et al., 2020a](#)).

4. JUNO

The Jiangmen Underground Neutrino Observatory (JUNO) is a large liquid scintillator detector currently under construction in an underground laboratory with a vertical overburden of about 650 m (roughly 1800 m water equivalent) in Jiangmen city in Southern China ([Adam et al., 2015](#)). JUNO is located 52.5 km from two nuclear power plants, a baseline optimized for JUNO’s primary goal, the determination of the neutrino mass ordering [28, 29]. To achieve this, JUNO requires a large target mass (20 kton) and excellent energy resolution, attributes also import in solar neutrino detection. The relatively shallow overburden limits the solar neutrino program, due to cosmogenic activity, but the low threshold and large detector mass may allow measurements of day/night effects, non-standard neutrino interactions af-

flecting the ^8B spectral shape, and ^7Be neutrinos, as well as providing a new solar determination of Δm_{12}^2 . JUNO’s reactor data will constrain Δm_{12}^2 to extremely high precision, allowing comparisons between the solar and reactor determinations of this parameter. To the extent that oscillation effects can be treated with greater confidence, the connections between solar neutrino flux measurements and the solar fusion reactions generating those fluxes will be sharpened ([An et al., 2016](#)).

5. DUNE

As recommended in the 2023 P5 report ([P52023, 2023](#)), the DUNE Collaboration seeks to construct two 10-kton fiducial liquid argon time projection chambers (LArTPC), deep underground in the Homestake mine in South Dakota ([Acciarri et al., 2016](#)). A third LArTPC will follow in Phase II, along with other upgrades. The newly developed Sanford Underground Research Facility (SURF) offers 4800-m water equivalent overburden, and will form the far site for a long baseline neutrino program utilizing a high-energy neutrino beam directed from Fermilab to SURF. The DUNE LArTPC detectors may be sensitive to measurements of the high-energy solar neutrinos via charged current interactions on argon, which offers good precision on the spectral shape and, thus, the potential to measure day/night effects ([Capozzi et al., 2019](#)). The solar neutrino program at DUNE is limited primarily by backgrounds (the detector is optimized for GeV-scale physics) and energy threshold. Multiple technologies are under consideration for a fourth detector module, with the motivation of expanding the physics program, including improved sensitivity to solar neutrinos. Use of underground argon, or alternative technologies such as organic or water-based scintillators, could preserve the long baseline neutrino sensitivity while also opening up a rich program of low-energy physics.

6. Future prospects

A “hybrid” detector that could utilize both Cherenkov and scintillation light simultaneously could achieve unprecedented levels of particle and event identification and, hence, background rejection ([Alonso et al., 2014](#)). The Cherenkov signature offers directional information, while the high light yield scintillation offers precision energy and vertex reconstruction. A full waveform analysis of detected light offers yet more information, based on the pulse shape of the scintillation, which is subject to species-dependent quenching effects, and the impact of the Cherenkov threshold. As a result, both the shape of the waveform and the ratio of the two signals will differ for different particle types. A full-scale detector utilizing novel scintillators along with fast and spectrally-sensitive

photon detectors, such as the proposed THEIA experiment (Askins *et al.*, 2020), or the Jinping detector (Beacom *et al.*, 2017) could achieve percent-level precision on the CNO neutrino flux, as well as improving precision across the suite of solar neutrino measurements, such as the ${}^8\text{B}$ spectral shape, and hep neutrino flux.

As experiments grow larger, and capabilities increase, there are also opportunities to apply detectors designed for other purposes to the detection of solar neutrinos. Noble liquid detectors built primarily for dark matter searches and long-baseline neutrino experiments have now reached the size and background levels to be sensitive to solar neutrinos (Aprile *et al.*, 2024). The detection of neutrinos from the high-energy tail of the ${}^8\text{B}$ spectrum via coherent elastic neutrino-nucleus scattering (“CE ν NS”) is particularly interesting because it is a purely neutral-current process sensitive to the total flux of active neutrinos (above the experimental threshold, about 8 MeV) from the Sun. Only the SNO experiment has had such a capability. Initial results reported in (Aprile *et al.*, 2024) are in agreement, with limited statistical precision. Solid state detectors built purposely for observing CE ν NS have observed neutrinos from stopped pion decay at the theoretically expected rate (Akimov *et al.*, 2017).

III. The ${}^1\text{H}(p, e^+\nu){}^2\text{H}$ reaction (S_{11})

The cross section for the initial reaction in the pp chain (Figure 1), $p + p \rightarrow d + e^+ + \nu_e$, is too small to be measured in the laboratory. It must be calculated from the standard theory of weak interactions.

A. Introduction and terminology

Near the Gamow peak energy $E \sim 6$ keV for temperatures characteristic of the solar core, the first and second derivatives of the astrophysical S factor at zero energy, $S'_{11}(0)$ and $S''_{11}(0)$, generate $\sim 7\%$ and $\sim 0.5\%$ corrections, respectively, in Taylor’s series expansion of $S_{11}(E)$ around $E = 0$. Higher derivative terms are neglected in this review since they only contribute at the 10^{-4} level. The recommended values for $S'_{11}(0)$ and $S''_{11}(0)$ will be discussed in Section III.E. Here we focus on $S_{11}(0)$.

At zero relative energy, $S_{11}(0)$ can be written as (Bahcall and May, 1968, 1969),

$$S_{11}(0) = 6\pi^2 m_p \alpha \ln 2 \frac{\bar{\Lambda}^2}{\gamma^3} \left(\frac{G_A}{G_V} \right)^2 \frac{f_{pp}^R}{(ft)_{0^+ \rightarrow 0^+}}, \quad (5)$$

where $\alpha = 1/137.04$ is the fine-structure constant, $m_p = 938.272$ MeV is the proton mass, $\gamma = (2\mu B_d)^{1/2} = 0.23161$ fm $^{-1}$ is the deuteron binding wave number, μ being the proton-neutron reduced mass and B_d the

deuteron binding energy, G_V and G_A are the Fermi vector and axial-vector weak coupling constants. Finally, f_{pp}^R is the phase-space factor for the pp reaction with radiative corrections, $(ft)_{0^+ \rightarrow 0^+}$ is the ft value for superallowed $0^+ \rightarrow 0^+$ transitions, and $\bar{\Lambda}$ is proportional to the transition matrix element connecting the pp and deuteron states.

B. Adopted parameters for this review

For the phase-space factor f_{pp}^R , we use the same value $f_{pp}^R = 0.144(1 \pm 0.001)$ as in SF II. It comes from the value without radiative corrections, $f_{pp} = 0.142$ (Bahcall and May, 1969), increased by 1.62% to take into account radiative corrections to the cross section (Kurylov *et al.*, 2003). The main source of uncertainty in f_{pp}^R arises from neglected diagrams in which the lepton exchanges a weak boson and a photon with different nucleons. These diagrams are estimated to modify f_{pp}^R by $\sim 0.1\%$, based on scaling the similar nucleus-dependent correction in superallowed β decay (Kurylov *et al.*, 2003). Direct computations of these diagrams were recommended in Solar Fusion II. Here we urge again that this computation be carried out.

For G_A/G_V , we use the value from the Particle Data Group (PDG) compilation (Workman *et al.*, 2022), 1.2754 ± 0.0013 , whose central value is larger than the 2008 PDG value, 1.2695 ± 0.0029 , used in SF II by 0.45% (or 1.9σ). Naively, this would lead to a $\sim 1\%$ increase in the central value of $S_{11}(0)$ according to Eq. (5). Its effect will be discussed below. For $(ft)_{0^+ \rightarrow 0^+}$, we take 3072.24 ± 1.85 s from the most updated comprehensive analysis of experimental rates with the radiative and Coulomb effects corrected (Hardy and Towner, 2020). This value is consistent with 3071.4 ± 0.8 s (Hardy and Towner, 2009) used in SF II with a larger error.

The dominant uncertainty in $S_{11}(0)$ comes from the normalized Gamow-Teller (GT) matrix element $\bar{\Lambda}$. Reducing this uncertainty has been the main focus of theoretical work since SF I. In SF I, $\bar{\Lambda}$ was written as $\bar{\Lambda} = \Lambda + \delta\Lambda$, i.e. the sum of the one- and two-body current matrix elements, Λ and $\delta\Lambda$, respectively, with their uncertainties estimated independently. In SF II, two major steps had contributed to reducing the uncertainty on $\bar{\Lambda}$. The first was a much deeper understanding of the correlation between the uncertainties in Λ and $\delta\Lambda$: the overall uncertainty in $\bar{\Lambda}$ could be described by a universal parameter that could be fixed by a single measurement. The study of Schiavilla *et al.* (1998) demonstrated this phenomenologically in the context of potential-model approaches, while later analysis via effective field theory (EFT) provided a more formal justification (Butler *et al.*, 2001; Park *et al.*, 2003). The second step was the use of the precisely known tritium β decay rate Γ_β^T to fix this universal parameter, as first proposed by Carlson *et al.*

(1991). This has been done in both potential models (Schiavilla *et al.*, 1998) and in the hybrid EFT approach (Park *et al.*, 2003) as explained below.

In SF II, $\bar{\Lambda}$ was determined with three approaches. The first one was the potential model approach. In the most elaborate calculation for the pp fusion process, a comparison of the results for five representative modern potentials designed to accurately reproduce nucleon-nucleon scattering data was carried out (Schiavilla *et al.*, 1998). After adjusting the unknown strength of the two-body exchange currents to reproduce Γ_β^T , the variation in $S_{11}(0)$ that otherwise would come from the choice of the phenomenological potential was largely removed. Predictions for five representative high-precision phenomenological potentials fell in a narrow interval $7.03 \lesssim \bar{\Lambda}^2 \lesssim 7.04$. There were additional uncertainties in the three-body potentials and three-body currents in Γ_β^T , of the order of $\sim 0.8\%$, and a 0.5% uncertainty due to effective range parameters for nucleon-nucleon scattering. Hence, the recommended $S_{11}(0)$ value from the potential model approach was $S_{11}(0) = 4.01(1 \pm 0.009) \times 10^{-25}$ MeV b.

The second and third approaches were both based on EFT. The second one was a hybrid EFT (EFT*), which used the current operators derived from EFT in conjunction with the initial and final state wave functions generated by a potential model (Park *et al.*, 2003). For pp fusion, the relevant two-body current contained only one unknown low-energy constant (LEC) \hat{d}^R which parameterized the contact axial coupling to two nucleons (Park *et al.*, 2003). A weakness of this approach was the mismatch between the operators and wave functions. However, it was argued that the mismatch only happened for short distance physics which could be absorbed by the LECs. Hence, when the ultraviolet cutoff was changed over a physically reasonable range, the residual cutoff dependence of physical observables provided a measure of the model dependence of the EFT* calculation. By combining the 0.8% error from changing the cutoff Λ_{NN} in the range of $500 - 800$ MeV and the $\sim 0.4\%$ higher order correction, obtained by multiplying the 1.8% contribution of the highest calculated order with the small expansion parameter $m_\pi/\Lambda_{NN} \sim 1/4$, Park *et al.* (2003) provided the value $S_{11}(0) = 4.01(1 \pm 0.009) \times 10^{-25}$ MeV b, in perfect agreement with the one obtained within the phenomenological approach.

The third study was performed with the pionless EFT ($\not\! EFT$) approach. It is a framework applicable to processes with the characteristic momentum p much smaller than the pion mass m_π , such that the pion field can be “integrated out” and becomes non-dynamical (Kaplan, 1996; Bedaque *et al.*, 1999; Chen *et al.*, 1999). In this approach, all nucleon-nucleon interactions and two-body currents are described by point-like contact interactions with a systematic expansion in powers of p/m_π . For all the deuteron weak breakup processes (e.g. νd and $\bar{\nu}d$ scattering) and their inverse processes, includ-

ing the pp fusion, only one two-body current (with coupling $L_{1,A}$) is needed up to next-to-next-to-leading order (N²LO) (Butler *et al.*, 2001). Therefore a single measurement will fix $L_{1,A}$ and the rates of all such processes. This feature is shared by the other approaches discussed above. The computation of $\bar{\Lambda}$ in $\not\! EFT$ was carried out to the second order in the p/m_π expansion by Kong and Ravndal (2001) and then to the fifth order by Butler and Chen (2001). Constraints on $L_{1,A}$ from two-nucleon systems (Butler *et al.*, 2002; Chen *et al.*, 2003) yielded $S_{11}(0) = 3.99(1 \pm 0.030) \times 10^{-25}$ MeV b.

Based on the consistent results of the above three approaches, SF II recommended $S_{11}(0) = 4.01(1 \pm 0.009) \times 10^{-25}$ MeV b. The SF I value $S'_{11}(0) = S_{11}(0)(11.2 \pm 0.1)$ MeV⁻¹ from Bahcall and May (1969) was not recomputed in SF II. However, with $S_{11}(0)$ reaching a 1% accuracy, new calculations of $S'_{11}(0)$ and $S''_{11}(0)$, together with the full pionful chiral EFT (χ EFT) computations to remove the unknown systematics of the hybrid EFT, were called for in SF II. These two challenges have been met in SF III and will be described below.

C. Experimental progress on muon capture of the deuteron

In SF II it was also recommended to carry out the experimental determination of the muon capture rate on deuteron, as proposed in the MuSun experiment (Kammel, 2003; Salvat and MuSun Collaboration, 2017). This quantity could be used to constrain $S_{11}(0)$ without the need to rely on the three-body calculation of Γ_β^T . The theoretical calculations for the muon capture rate on deuteron have been carried out in $\not\! EFT$ (Chen *et al.*, 2005) (it is possible to impose neutron energy cut to isolate the low-energy neutron events so $\not\! EFT$ is applicable (Kammel, 2003)), chiral hybrid EFT* (Ando *et al.*, 2002), phenomenological potential model (Marcucci *et al.*, 2011), and recently in χ EFT (Marcucci *et al.*, 2012, 2018; Acharya *et al.*, 2018; Ceccarelli *et al.*, 2023; Bonilla *et al.*, 2023; Gnech *et al.*, 2024). However, the MuSun result is yet to be released. On the other hand, we should notice that the muon capture processes happen at a rather large momentum transfer compared to pp fusion. The momentum transfer dependence of the single nucleon axial coupling constant $g_A(q^2)$, with $g_A(q^2 = 0) \equiv g_A \equiv G_A/G_V$, has been recently studied by Hill *et al.* (2018), who have provided an experimental determination for the axial charge radius, given by $r_A^2 = 0.46(16)$ fm², where r_A is defined by the relation $g_A(q^2) = g_A(1 - r_A^2 q^2/6)$ for small q^2 . The $\sim 30\%$ uncertainty on r_A^2 has an impact on the ability of the MuSun experiment alone to directly constrain $S_{11}(0)$ (Acharya *et al.*, 2018; Bonilla *et al.*, 2023; Ceccarelli *et al.*, 2023; Gnech *et al.*, 2024).

In the next few years, lattice Quantum Chromodynamics (QCD) calculations of $g_A(q^2)$ are expected to reduce

the r_A uncertainty by a factor of two or more (see Meyer (2023) for the most recent review of the lattice results). Currently, lattice QCD results are consistent with Hill *et al.* (2018) on r_A , often with comparable uncertainties. At larger q^2 ($|q^2| \gtrsim 0.25$ GeV 2), there is a growing tension between lattice QCD predictions of $g_A(q^2)$ and the phenomenological determination from older neutrino-deuterium bubble chamber data (Meyer *et al.*, 2016) with the lattice QCD results yielding a 30% larger neutrino-nucleon cross section over a large range of q^2 (Meyer *et al.*, 2022).

In any case, it is evident that accurate experimental determinations of muon capture rates on the deuteron and other light nuclei, which can be addressed theoretically *ab initio* approaches, represent fundamental tests for the theoretical approaches themselves, either within χ EFT or π EFT, and might be able to provide the necessary experimental information to fix the unknown parameters of the theory.

D. Progress in $S_{11}(0)$ calculations since SF II

Below we summarize the $S_{11}(0)$ calculations performed after SF II using different approaches.

1. χ EFT

The pioneering work of Marcucci *et al.* (2013) used the next-to-next-to-next-to-leading-order (N^3LO) chiral two-nucleon potential (Entem and Machleidt, 2003; Machleidt and Entem, 2011) augmented with higher order ($\mathcal{O}(\alpha^2)$) two-photon and vacuum-polarization electromagnetic interactions. These $\mathcal{O}(\alpha^2)$ corrections reduced $S_{11}(0)$ by $\sim 0.8\%$ mainly due to the vacuum-polarization-induced pp wave function distortion. This was consistent with the 0.84% first found in the potential model calculation of Schiavilla *et al.* (1998). This correction was also included in the EFT* calculation of (Park *et al.*, 2003). The relevant LECs were fitted to reproduce the $A = 3$ binding energies, magnetic moments, and GT matrix element in Γ_β^T to obtain $S_{11}(0) = 4.030(1 \pm 0.006) \times 10^{-25}$ MeV b, with the p -wave initial state contributing at $\sim 1\%$, which was about the accuracy level of the calculation (Marcucci *et al.*, 2013). However, using π EFT at next-to-leading-order (NLO), Acharya *et al.* (2019) later found that p -wave only contributed at the order of 10^{-30} MeV b. Re-examining the computer programs, Marcucci *et al.* (2013) found an error in the determination of one of the p -wave reduced matrix elements (associated with the longitudinal multipole operator). Consequently, Marcucci *et al.* (2019) reported in the Erratum that $S_{11} = 4.008(1 \pm 0.005) \times 10^{-25}$ MeV b.

In more recent work, Acharya *et al.* (2016) used chiral interactions and consistent currents up to N^3LO (called

next-to-next-to-leading-order (N^2LO) in the original literature since the second order vanished) and developed a robust procedure for the error quantification. In particular, they analyzed a family of 42 interactions (Carlsson *et al.*, 2016) with 7 different cutoff values from 450 to 600 MeV. The 26 LECs were fitted to 6 different pools of input data including NN and πN scatterings, as well as the binding energies and charge radii of 3H and 3He , the quadrupole moment of 2H , and Γ_β^T . This thorough study yielded $S_{11}(0) = 4.047(1^{+0.006}_{-0.008}) \times 10^{-25}$ MeV b.

In both of the χ EFT calculations mentioned above, a widely used relation first proposed by Gazit *et al.* (2009) that linked the two-body axial current LEC \hat{d}_R with the LEC c_D from the πNN vertex was employed. However, later Marcucci *et al.* (2018) found that there was a factor $-1/4$ missing in this \hat{d}_R - c_D relation which was then acknowledged in the Erratum of Gazit *et al.* (2019). Fortunately, this error was unimportant in muon capture on deuteron (Marcucci *et al.*, 2018) and it only affected $S_{11}(0)$ at the 0.1% level (Acharya *et al.*, 2023).

In the most recent and comprehensive χ EFT study, Acharya *et al.* (2023) compared the above calculations of Marcucci *et al.* (2019) and Acharya *et al.* (2016) in detail. In addition to the 0.1% increase of $S_{11}(0)$ from using the correct \hat{d}_R - c_D relation, updating the input parameters to their most recent values increased $S_{11}(0)$ by $\sim 1\%$ in both calculations, mainly due to the 0.45% increase of G_A/G_V from its SF II value mentioned above. Furthermore, Marcucci *et al.* (2019) received a $\sim 1\%$ increase from removing the truncation error of the basis functions which effectively cut off the long distance part of the wave functions, as Acharya *et al.* (2017) advocated. After these corrections, the $\sim 1\%$ difference between Marcucci *et al.* (2019) and Acharya *et al.* (2016) was reconciled and the combined result was found to be (Acharya *et al.*, 2023)

$$S_{11}(0) = 4.100(1 \pm 0.007) \times 10^{-25} \text{ MeV b .} \quad (6)$$

Acharya *et al.* (2023) obtained consistent values of $S_{11}(0)$ using four different χ EFT models for the nuclear interaction. This value is also in agreement with the result obtained in π EFT by De-Leon and Gazit (2022) (see below). In estimating the order-by-order convergence, however, Acharya *et al.* (2023) used one of the χ EFT models, which is able to nicely reproduce the deuteron properties already at leading order. Therefore, the χ EFT error of Eq. (6) is likely to be an underestimate, warranting the enlarged error advocated in Section III.D.4.

Finally, we would like to mention the work of Liu *et al.* (2022a), where the power counting of the χ EFT weak current operator involved in the pp reaction is revisited using renormalization group (RG) invariance as the guideline. In particular, it is argued that the contact two-body axial current proportional to the \hat{d}_R LEC must appear one order lower than assessed by naive dimensional analysis. Then it can be shown that RG in-

variance is fulfilled at N²LO. However, the estimate for $\bar{\Lambda}$ obtained by Liu *et al.* (2022a) does not use a value for \hat{d}_R obtained by fitting Γ_β^T , but extracted in order to match the value of $\bar{\Lambda}$ obtained in SF II. Therefore, we will not consider the work of Liu *et al.* (2022a) in the present $S_{11}(0)$ evaluation.

2. $\not\!\! EFT$

The universal two-body current coupling $L_{1,A}$ was determined using Γ_β^T in $\not\!\! EFT$ for the first time by De-Leon *et al.* (2019). This calculation was carried out up to NLO using the dibaryon formulation of Beane and Savage (2001), which partially resummed higher order effective range contributions to improve the convergence. Then, this result was used by De-Leon and Gazit (2022), with updated input parameters, to obtain

$$S_{11}(0) = 4.12(1 \pm 0.015) \times 10^{-25} \text{ MeV b}, \quad (7)$$

where the $\mathcal{O}(\alpha^2)$ electromagnetic correction was not calculated but assumed to be the same as the potential model value, 0.84% (Schiavilla *et al.*, 1998). While this number was not model independent, the model dependence was believed to be well below the assigned 1.5% error.

The small error assigned to this NLO result has been justified by drawing an analogy from the corresponding electromagnetic processes. Using the same approach, the $np \rightarrow d\gamma$ matrix element at threshold was predicted at NLO within 0.5% to the experimental value after the electromagnetic two-body current L_1 was fit to the magnetic moments of ³He and ³H. This indicates that the contribution of the three-body current at N²LO is small in this electromagnetic case. The weak sector is shown to follow the same operator structure and hence provides support for the calculation procedure and the uncertainty estimate.

It is important to note that the reason that the three-body current is an N²LO effect in $\not\!\! EFT$ is related to the non-trivial renormalization of the non-derivative three-body contact interaction, which shows up at LO to absorb the cutoff dependence of Feynman diagrams. The subleading two-derivative three-body contact interaction is expected to show up at N²LO. This interaction, combined with the one-body current, renormalizes the three-body current. Hence, the three-body current should also appear at the same order, N²LO. However, if the non-derivative three-body contact interaction were counted as higher order, such as N³LO, as in χEFT (because cutoff independence is not strictly enforced order by order in χEFT), then the three-body current would contribute at much higher order. Although the χEFT power counting indeed yields good convergence in the expansions, it is unsatisfactory that one can not remove the cutoff dependence at each order of the expansion. In addition,

the uncertainty estimate of the χEFT still lacks a broad inspection of the specific nuclear χEFT potential implementation. As a consequence, we consider the $\not\!\! EFT$ 1.5% error a better estimate for the $S_{11}(0)$ theoretical uncertainty.

3. Lattice QCD and Lattice EFT

Ideally lattice QCD would provide a first principles prediction of the pp fusion rate and the GT matrix element of Γ_β^T , for both pure QCD and with QED effects incorporated. However, such calculations are quite challenging and not yet available at the required precision. A proof-of-principle calculation was carried out by Savage *et al.* (2017) using a background field method to determine both the pp fusion GT matrix element and Γ_β^T . This exploratory calculation utilized a single pion mass at the SU(3)-flavor-symmetric point with $m_\pi \approx 806$ MeV, a single and relatively coarse lattice spacing of $a \approx 0.145$ fm, and a single volume. The calculation was performed under the assumption that the two- and three-nucleon systems were deeply bound. Without this assumption, matrix elements computed in the finite volume can be significantly different from those in infinite volume, due to Lellouch-Lüscher factors (Lellouch and Lüscher, 2001) that lead to power-law finite volume corrections. These can range from the few-percent level to O(1) (Briceño and Davoudi, 2013; Briceño *et al.*, 2015).

More recent lattice QCD calculations have found that two-nucleon systems at heavy pion masses are in fact not bound (Iritani *et al.*, 2016; Francis *et al.*, 2019; Hörz *et al.*, 2021; Amarasinghe *et al.*, 2021). These efforts, which employ interpolating operators more sophisticated than those of Savage *et al.* (2017), suggest there could be large systematic uncertainties affecting the conclusions of Savage *et al.* (2017) stemming from misidentification of the spectrum and inaccurate Lellouch-Lüscher factors. In addition, Green *et al.* (2021) found that the two-baryon spectrum may be particularly sensitive to discretization effects, which would also have important implications for the continuum extrapolation of the matrix elements. These issues are discussed in some detail in Tews *et al.* (2022). Finally, results with pion masses $m_\pi \lesssim 300$ MeV are needed for accurate extrapolation to the physical pion mass.

Given these unresolved systematic issues, the result of Savage *et al.* (2017) is not included in the present $S_{11}(0)$ evaluation – even though the extracted value $S_{11}(0) = 4.07(1 \pm 0.008) \times 10^{-25}$ MeV b is consistent with our recommended range quoted in Section III.D.4.

The lattice EFT computation of Rupak and Ravi (2015) performed the pp fusion calculation by implementing $\not\!\! EFT$ on a spacetime lattice. The purpose of this leading-order study was to demonstrate that lattice EFT could reproduce the infinite volume and continuum re-

sult of $\#$ EFT such that it could be applied to various reactions of astrophysical interest in the future. Therefore, for $S_{11}(0)$, this result is considered as a subset of the $\#$ EFT calculation.

4. Final Recommendation of $S_{11}(0)$

The above discussions show that determinations of $S_{11}(0)$ from χ EFT in Eq.(6), $\#$ EFT in Eq.(7), and lattice QCD (although with unquantified systematics) are all consistent with each other. Furthermore, these values are also consistent with the recommended value of SF II, provided the central value is increased by 0.9%⁵ to $S_{11}(0) = 4.05(1 \pm 0.009) \times 10^{-25}$ MeV b to account for the G_A/G_V update. Averaging this value with the χ EFT value in Eq.(6) and the $\#$ EFT value in Eq.(7) yields $S_{11}(0) = 4.09(1 \pm 0.005) \times 10^{-25}$ MeV b with χ^2 per degree of freedom to be 0.9. This shows that the χ EFT, $\#$ EFT and SF II estimates of $S_{11}(0)$ are all mutually consistent.

In addition, we would like to advocate adding an additional correlated error to account for any input that would tend to move all results in a coordinated way. For example, from SF II to SF III, we experienced the $\sim 1\%$ shift due to the update of the PDG value of G_A/G_V . It is not inconceivable that G_A/G_V or other input parameters or physics could change again by similar amounts: the large PDG inflation factor of 2.7 reflects the tension that continues to exist among G_A/G_V measurements (Workman *et al.*, 2022). Therefore, assigning an additional 1% correlated error seems reasonable. This is also in line with the subtleties discussed in Secs.III.D.1 and III.D.2 which call for an enlarged error. Therefore, our final recommended value for $S_{11}(0)$ is

$$S_{11}(0) = 4.09(1 \pm 0.015) \times 10^{-25} \text{ MeV b}, \quad (8)$$

where we have added the correlated and uncorrelated errors linearly to be conservative.

E. Progress in $S'_{11}(0)$ and $S''_{11}(0)$

Using $\#$ EFT, Chen *et al.* (2013) computed $S_{11}(E)$ analytically with all partial waves included up to N²LO. The Fermi matrix element only contributed at the 10^{-4} level and was neglected compared with the GT matrix element. The energy dependence of the phase factor f_{pp}^R of Eq. (5) was the dominant effect in $S'_{11}(0)$ and $S''_{11}(0)$, the energy dependence of pp scattering was subdominant,

while the $L_{1,A}$ contribution was much less important in these derivatives than in $S_{11}(0)$. Therefore, these derivatives could be predicted more reliably than $S_{11}(0)$. Furthermore, the derivatives were computed analytically and were free from errors of fitting $S_{11}(E)$ to a polynomial. The result was $S'_{11}(0)/S_{11}(0) = (11.3 \pm 0.1)$ MeV⁻¹ and $S''_{11}(0)/S_{11}(0) = (170 \pm 2)$ MeV⁻².

In χ EFT, $S_{11}(E)$ for $E < 100$ keV was fit to polynomials of E (Marcucci *et al.*, 2019). Depending on using a quadratic or quartic fit, $S'_{11}(0)/S_{11}(0)$ changed from 12.23 to 10.82 MeV⁻¹ and $S''_{11}(0)/S_{11}(0)$ changed from 178.4 to 317.4 MeV⁻². Acharya *et al.* (2016) used cubic fit and $E < 30$ keV to obtain $S'_{11}(0)/S_{11}(0) = 10.84(2)$ MeV⁻¹ and $S''_{11}(0)/S_{11}(0) = 317.8(13)$ MeV⁻². Recently Acharya *et al.* (2023) repeated the calculation of Marcucci *et al.* (2019) with the same energy range and cubic fit as Acharya *et al.* (2016), obtaining a consistent result with Acharya *et al.* (2016): $S'_{11}(0)/S_{11}(0) = 10.83$ MeV⁻¹ and $S''_{11}(0)/S_{11}(0) = 313.72$ MeV⁻². We will take these as the recommended values from χ EFT.

Although the face values of $S'_{11}(0)/S_{11}(0)$ and $S''_{11}(0)/S_{11}(0)$ from $\#$ EFT and χ EFT look quite different, we would like to remark that they actually agree on $S_{11}(E)/S_{11}(0)$ better than 0.1% below the ~ 6 keV Gamow peak. For massive stars with central temperatures ~ 15 keV, the agreement is better than 0.8% to second order in the derivatives and 0.5% if $S'''_{11}(0)/S_{11}(0) = -5382$ MeV⁻³ is included in the χ EFT result. Hence, we take the average of the $\#$ EFT and χ EFT results as the recommended value:

$$\begin{aligned} S'_{11}(0) &= S_{11}(0)(11.0 \pm 0.2)\text{MeV}^{-1}, \\ S''_{11}(0) &= S_{11}(0)(242 \pm 72)\text{MeV}^{-2}. \end{aligned} \quad (9)$$

Finally, we comment on the work of Gaspard *et al.* (2019) which performed a study of $S_{11}(E)$ in a wide energy range. The main focus of this work was to perform a proper energy-dependence analysis of the pp process, in order to reliably extract $S'_{11}(0)/S_{11}(0)$ and $S''_{11}(0)/S_{11}(0)$. However, the calculation was performed within a phenomenological approach, using a quite simplified model for the nuclear currents (i.e. not including two-body currents, which are well known to be significant) and structure (i.e. neglecting the d -wave components in the deuteron wave function). They found results for $S'_{11}(0)/S_{11}(0)$ and $S''_{11}(0)/S_{11}(0)$ compatible with those of Eq. (9). However, since these values were not obtained with state of the art calculations, they have not been considered in the determination of $S'_{11}(0)/S_{11}(0)$ and $S''_{11}(0)/S_{11}(0)$.

⁵ To know the precise shift requires an explicit calculation. However, if the shift is within the range of 0.8-1.0 %, the averaged value and χ^2 remain the same within the significant digits.

IV. The $^2\text{H}(p,\gamma)^3\text{He}$ reaction (S_{12})

A. Introduction

The $^2\text{H}(p,\gamma)^3\text{He}$ reaction is the second step in the pp chain (Figure 1). Compared with the reactions mediated by the weak interaction, this reaction occurs much more rapidly. Consequently, on the time scales relevant to solar energy generation, deuterium is effectively converted to ^3He instantaneously and thus it is only sensitive to the Q value of the reaction and not the uncertainties of the rate.

However, the $^2\text{H}(p,\gamma)^3\text{He}$ reaction plays an important role in the development of protostars, because the onset of deuterium burning slows down the protostars contraction and heating, increasing their lifespan. Accurate knowledge of the $^2\text{H}(p,\gamma)^3\text{He}$ reaction rate, particularly within the few keV range corresponding to the Gamow peak in protostars, is vital for modeling protostellar evolution effectively (Stahler, 1988).

Another astrophysical scenario where the $^2\text{H}(p,\gamma)^3\text{He}$ reaction plays a key role is Big Bang Nucleosynthesis (BBN), responsible for the production of light elements during the first few minutes of the Universe. Among these elements, deuterium is an excellent indicator of cosmological parameters because its primordial abundance is the most sensitive to the baryon density and critically depends on the radiation density of the early Universe, see for example a review of BBN (Cyburt *et al.*, 2016). The reactions involved in the synthesis of deuterium are: production via the well known $p(n,\gamma)^2\text{H}$ process and destruction via the $^2\text{H}(^2\text{H},n)^3\text{He}$, $^2\text{H}(^2\text{H},p)^3\text{H}$ and $^2\text{H}(p,\gamma)^3\text{He}$ reactions (Fields *et al.*, 2020).

Since the comprehensive review performed in SF II, there have been both new measurements as well as advances in the theoretical and phenomenological analysis of the $^2\text{H}(p,\gamma)^3\text{He}$ reaction. The new experimental results have been determined with accelerator-based measurements of the cross section (Mossa *et al.*, 2020b; Tišma *et al.*, 2019; Turkat *et al.*, 2021), as well as plasma-based, inertial confinement fusion measurements (Mohamed *et al.*, 2022; Zylstra *et al.*, 2020) and are reviewed in Section IV.B below. On the theoretical side, there have been advances in *ab initio* calculations, where nucleons are the fundamental degrees of freedom interacting among themselves and with the external electromagnetic probe, see for example (Ekström *et al.*, 2023). Finally, Bayesian analysis methods have been used to model the energy dependence of the S-factor, starting from the *ab initio* predictions and applying a polynomial approximation to it.

We review these updates and provide recommended values and uncertainties of the S-factor over the energy range of interest for Solar fusion, based upon a Bayesian averaging of various models.

B. Data sets used for the present review

The $^2\text{H}(p,\gamma)^3\text{He}$ reaction has a Q value of 5.5 MeV and proceeds through the radiative capture of a proton on deuterium. Different experimental approaches were followed to measure its cross section. Tišma *et al.* (2019) irradiated deuterated titanium targets with a proton beam and detected the γ -rays with two high-purity germanium (HPGe) detectors placed at different angles. The final S-factor is provided at four energies in the 47 - 210 keV range, with approximately 15% statistical uncertainty. More recently, the LUNA Collaboration performed a measurement in the energy range of 32 – 263 keV at the underground in the Gran Sasso Laboratories, exploiting the six orders of magnitude suppression of the cosmic radiation background (Cavanna and Prati, 2018; Ferraro *et al.*, 2021). A windowless deuterium gas target was used and the γ -rays emitted by the $^2\text{H}(p,\gamma)^3\text{He}$ reaction were detected by a large HPGe detector at 90° with respect to the beam axis. Great care was taken to minimize all sources of systematic uncertainties in the S-factor overall uncertainty at the 3% level (Mossa *et al.*, 2020a,b). These new results provided stringent constraints on cosmological parameters obtained by comparing the precise primordial deuterium abundance predictions of the standard BBN model with astronomical observations (Cooke *et al.*, 2018). A deeper discussion of the LUNA results and their implications can be found in (Moscoso *et al.*, 2021; Pisanti *et al.*, 2021; Pitrou *et al.*, 2021; Yeh *et al.*, 2021)⁶

Finally, a new measurement was performed at the Helmholtz-Zentrum Dresden-Rossendorf in the 300–1000 keV energy range using implanted deuterium targets on tantalum backings and two HPGe detectors (Turkat *et al.*, 2021). The resulting S-factors show $\approx 1 - 2\sigma$ tension with the analysis of the LUNA results extrapolated to $E \gtrsim 300$ keV (Mossa *et al.*, 2020b). However, they are affected by large systematic uncertainties. A new measurement of the $^2\text{H}(p,\gamma)^3\text{He}$ reaction is planned at the Felsenkeller laboratory in Germany (Bemmerer *et al.*, 2019).

In addition to the new accelerator-based results, two recent sets of measurements (Mohamed *et al.*, 2022; Zylstra *et al.*, 2020) have also been performed using the inertial confinement plasma-based platform (Gatu Johnson *et al.*, 2023, 2017; Mohamed *et al.*, 2022), which has recently begun to be exploited for this type of work (e.g., Casey *et al.*, 2017; Zylstra *et al.*, 2016). Both measurements, which were performed at the OMEGA laser

⁶ After the closing of the present review, the Mossa *et al.* (2020b) data were re-analyzed to extract the experimental $^2\text{H}(p,\gamma)^3\text{He}$ γ -ray angular distribution (Stöckel *et al.*, 2024). The data confirm the ab initio angular distributions used in the original analysis by Mossa *et al.* (2020b).

facility (see section XIV.D), used laser-driven implosions of spherical plastic-shell capsules filled with H_2D_2 gas, and measured the emitted γ -rays using a gas-Cherenkov-detector (Herrmann *et al.*, 2014) that was calibrated applying the technique described in Zylstra *et al.* (2019). The initial experiment (Zylstra *et al.*, 2020) obtained good statistics at an energy of 16 keV by making several repeated measurements, with a final statistical uncertainty of 6% and a systematic one of 17% (dominated by uncertainty in the absolute calibration of the detector). The second experiment (Mohamed *et al.*, 2022) obtained data at three different energies in the 17–37 keV region, with comparable systematic uncertainty, but with larger statistical uncertainty due to fewer repeated measurements. The results obtained on this unique platform agree within error bars with the accelerator-based measurements.

C. Theoretical studies

Nuclear reactions of astrophysical interest in general, and the ${}^2\text{H}(p,\gamma){}^3\text{He}$ in particular, are of great importance in nuclear theory because the available experimental data can be used to test the adopted theoretical framework. The ${}^2\text{H}(p,\gamma){}^3\text{He}$ reaction has the great advantage of involving only $A \leq 3$ nuclear systems, and can be addressed with a microscopic *ab initio* study. This means that the nuclear systems involved in the process are viewed as made up of A nucleons, interacting among themselves and with the external electromagnetic probes. Within such an approach, the following ingredients are essential for the calculation: (i) realistic models for the nuclear interactions and currents, possibly rooted in Quantum Chromodynamics (QCD); (ii) a numerical technique able to solve the A -body bound and scattering state problem, including the Coulomb interaction without approximation. Such a technique is usually referred to as an *ab initio* method. The agreement (or disagreement) between *ab initio* theoretical predictions and experimental data represents a validation (or indicates the necessity of improvement) mostly for ingredient (i), i.e. are the models of the nuclear interactions and currents accurate within the precision specified by the *ab initio* method? This is why few-nucleon reactions can be used as an “ideal” laboratory, where the *ab initio* framework can be stringently tested in systems under more theoretical control.

The most recent *ab initio* calculation of the ${}^2\text{H}(p,\gamma){}^3\text{He}$ reaction is that of Marcucci *et al.* (2016). Here the pair-correlated Hyperspherical Harmonics *ab initio* method was used to calculate the $A = 3$ initial scattering and final bound state wave functions (see Kievsky *et al.* (2008) and Marcucci *et al.* (2020) for details). The nuclear interaction model adopted to describe the $A = 3$ nuclear state in Marcucci *et al.* (2016) consists of a two-nucleon term, the Argonne v_{18} (AV18) potential (Wiringa *et al.*, 1995),

augmented by a three-nucleon contribution, the Urbana IX (UIX) potential (Pudliner *et al.*, 1995). The AV18 potential can reproduce the large two-nucleon database with a $\chi^2/\text{datum} \sim 1$ (Wiringa *et al.*, 1995), while the combination AV18/UIX can describe quite accurately the properties of ${}^3\text{He}$, the spectra of light p-shell nuclei (Pieper and Wiringa, 2001), and $p - d$ scattering observables (see for instance Wood *et al.* (2002)). The electromagnetic current operator used in Marcucci *et al.* (2016) includes, in addition to the non-relativistic one-body operator, two- and three-body terms required by gauge invariance in a system of interacting particles. These terms were constructed in Marcucci *et al.* (2005). The model was then tested against various electromagnetic observables, to access the quality of its predictions. As a potential model, however, there is no systematic procedure for assigning uncertainties for observables whose values are unknown.

The results of Marcucci *et al.* (2016) have been found to be about 10% higher than the experimental data of Mossa *et al.* (2020b). Given the lack of a procedure for quantifying errors in calculations based on phenomenological interactions and currents, it is difficult to access the significance of this discrepancy. This leads us to make several recommendations to the theory community. First, the ${}^2\text{H}(p,\gamma){}^3\text{He}$ reaction should be studied within the framework of chiral effective field theory, which has reached a degree of accuracy and predictive power comparable to potential-based phenomenology. This approach is formally *ab initio* and rooted in QCD, and as an operator expansion will provide an estimate of the theoretical uncertainty. Work along this line is currently underway. Second, the angular distribution of the ${}^2\text{H}(p,\gamma){}^3\text{He}$ capture reaction, and possibly also polarization observables, should be both calculated and measured. This would be valuable even if measurements were limited to higher energies, where they are less difficult. Such measurements would provide a further test of the predictive power of the theory.

D. Phenomenological and Bayesian Analyses

In the energy range of astrophysical interest, the ${}^2\text{H}(p,\gamma){}^3\text{He}$ reaction does not have any resonance or coupled channels that can give rise to non-trivial energy dependence, and the S-factor can be modelled by a low-order polynomial in energy (Fowler *et al.*, 1967). Moreover, the recent results from LUNA (Mossa *et al.*, 2020b), combined with previous measurements, place stringent constraints on the cross section, reducing the uncertainty over the range of interest to Solar fusion and BBN (Casella *et al.*, 2002; Mossa *et al.*, 2020b; Pisanti *et al.*, 2021; Pitrou *et al.*, 2021; Yeh *et al.*, 2021).

A significant change since SF II has been the wide adoption of Bayesian analysis methods to evaluate ther-

TABLE III Logarithm of the Gaussian approximation to the Bayes factor, and corresponding weight, Eq. (B4), for each model under investigation.

Model	logGBF	weight
<i>ab initio</i> , $aS_{\text{nuc}}(E)$	1478.7	0.585
<i>ab initio</i> , $aS_{\text{nuc}}(E) + b$	1478.3	0.414
3 rd order polynomial	1470.2	1.2×10^{-4}
4 th order polynomial	1469.2	4.6×10^{-5}
5 th order polynomial	1468.2	1.6×10^{-5}
6 th order polynomial	1464.1	2.9×10^{-7}

monuclear reaction rates and provide more rigorous uncertainty estimates in a statistical sense (Iliadis *et al.*, 2016). This approach has been followed in the analysis presented here, which closely follow work on the ${}^2\text{H}(p,\gamma){}^3\text{He}$ reaction reported in Moscoso *et al.* (2021), and references therein. The mathematical details of the Bayesian method will be found in Appendix B: Here we focus on the application to ${}^2\text{H}(p,\gamma){}^3\text{He}$ by the authors cited above.

Both the *ab initio* prediction from Marcucci *et al.* (2005) as well as a third-order polynomial in the energy were used to constrain the data, where we define an n^{th} -order polynomial as

$$S_n(E; \lambda) = \sum_{i=0}^n \lambda_i E^i. \quad (10)$$

As noted in the literature, the energy dependence of Marcucci *et al.* (2005) is more reliable than the absolute normalization. Furthermore, the updated prediction in Marcucci *et al.* (2016), as noted above, is $\approx 10\%$ larger than that of Marcucci *et al.* (2005). Therefore, Moscoso *et al.* (2021) modeled the prediction of Marcucci *et al.* (2005), denoted $S_{\text{nuc}}(E)$, as

$$S(E; \lambda) = aS_{\text{nuc}}(E) + b, \quad (11)$$

where a and b are an unknown scale factor and offset to be determined in the analysis. While $S_{\text{nuc}}(E)$ was not determined with theoretical uncertainty, model uncertainty is introduced through the parameters a and b . The resulting mean values and uncertainties of $S(E; \lambda)$ determined from Eq. (11) were found to be comparable to those with the third-order polynomial.

E. Summary and Recommendations

In this review, we take the further step of applying Bayes Model Averaging to a number of reasonable models that describe the data, which systematically captures additional theoretical “model selection uncertainty”. Details of the Bayesian analysis and model averaging procedure are presented in Appendix B. The models we explore

include Eq. (11) with both a and b , as done in Moscoso *et al.* (2021), as well as setting $b = 0$, and then we consider polynomials in energy of order $n = 3, 4, 5, 6$.⁷

With such a Bayes model averaging, we can quantitatively compare and contrast the polynomial parameterizations of ${}^2\text{H}(p,\gamma){}^3\text{He}$ along with *ab initio* results predicted by Marcucci *et al.* (2016, 2005), including the extra variance that arises from this set of reasonable models, e.g., using polynomials of different order, as well as the phenomenological model of Eq. (11). Results are listed in Table III. The first column lists the model (i.e., polynomial or *ab initio*), the second column the natural logarithm of the Gaussian approximation to the BF, and the third column the corresponding weight in the model averaging. From this model averaging, the resulting prediction for the ${}^2\text{H}(p,\gamma){}^3\text{He}$ S-factor at a few representative energies is provided in Table IV, where the first uncertainty arises from the first term in Eq. (B3) and the second uncertainty is from the second term, which we denote as *model selection uncertainty*. The resulting Bayes model average prediction of the S-factor over the entire kinematic range considered is depicted in Fig. 5, with the gray band representing the 68% coverage probability. For this particular reaction, it is interesting to note that the Bayesian analysis strongly favors the phenomenological models of Eq. (11) over the polynomial approximations. One reason for this might be that the scaled *ab initio* models have only one or two free parameters, and the energy dependence of $S_{\text{nuc}}(E)$ given in Marcucci *et al.* (2016, 2005) is sufficient to accurately describe the various data sets. On the other hand, polynomial approximations are disfavored as they require more parameters to capture the energy dependence. Also of note, in the model of Eq. (11) with $b = 0$ and using Marcucci *et al.* (2016) for $S_{\text{nuc}}(E)$, the scale factor is given by

$$a = 0.921(19), \quad (12)$$

indicating that the prediction in Marcucci *et al.* (2016) overestimates the ${}^2\text{H}(p,\gamma){}^3\text{He}$ data by 7.9%, consistent with expectations noted above. In comparison, the third order polynomial fit predicts values of $S(E)$ that are 1σ higher at $E = 0$ and $\frac{2}{3}\sigma$ lower at $E = 91$ keV as compared to those in Table IV.

The analysis in Table III and Fig. 5 can be reproduced with the code linked in Table IV. The code also can provide a prediction of the mean value and uncertainty of $S_{12}(E)$ at any energy over the same kinematic range.

⁷ The implementation of the Bayesian analysis presented here is different than that of Moscoso *et al.* (2021) but it was verified to produce the same results when the same implementation of the systematic uncertainties was made.

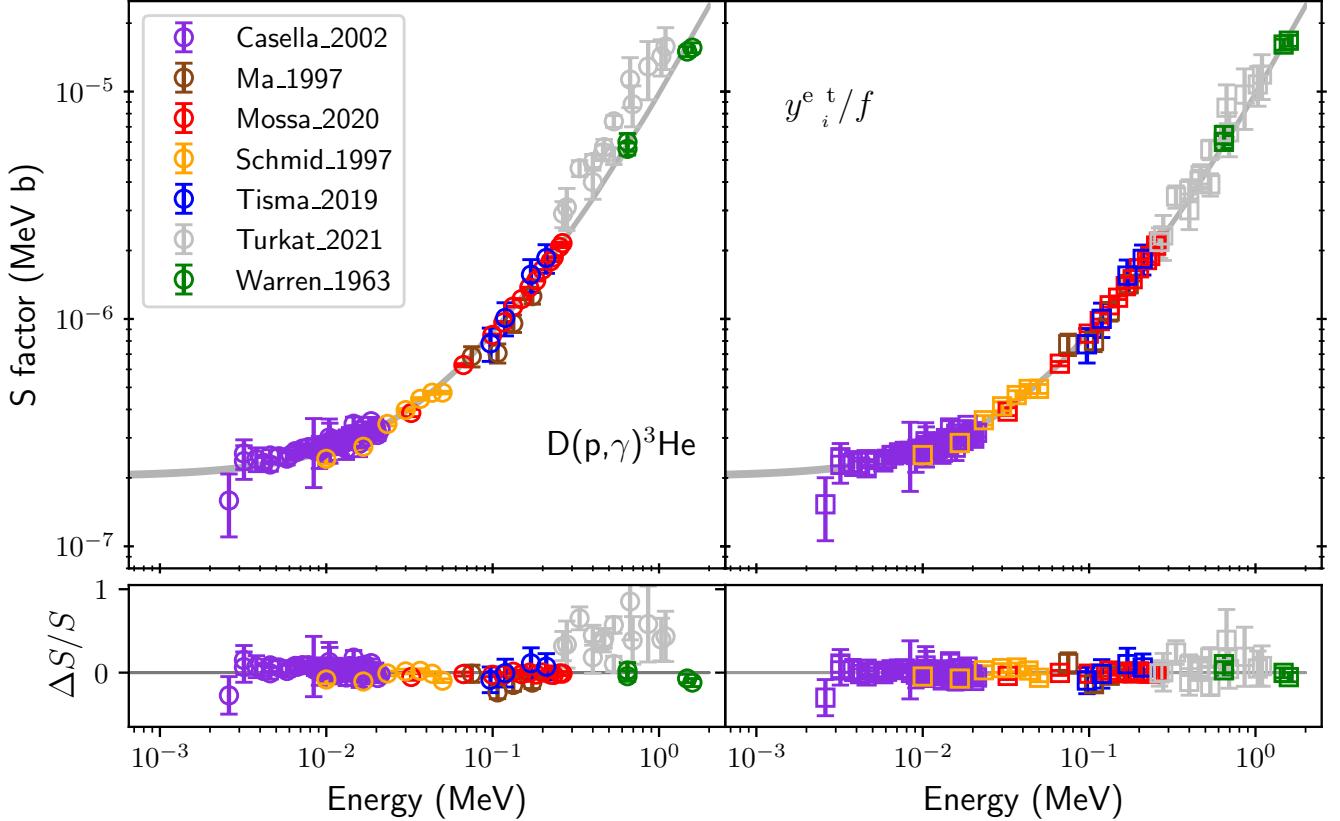


FIG. 5 S-factor determined via Bayes model averaging (BMA) of the ${}^2\text{H}(p,\gamma){}^3\text{He}$ data sets (Casella *et al.*, 2002; Ma *et al.*, 1997; Mossa *et al.*, 2020b; Schmid *et al.*, 1997; Tišma *et al.*, 2019; Turkat *et al.*, 2021; Warren *et al.*, 1963) analysed as described in the text. The left panel shows the original data with statistical uncertainties only. The right panel shows the data after normalizing by BMA systematic uncertainty posteriors of f_D^{-1} for each data set, and the uncertainties correspond to the statistical, systematic and extrinsic uncertainties all added in quadrature. In most cases, the quoted statistical uncertainty is the dominant uncertainty. The residual is defined as $\Delta S/S = (S_{\text{data}} - S_{\text{fit}}^{\text{BMA}})/S_{\text{fit}}^{\text{BMA}}$ and is plotted for the original data with statistical uncertainty only on the left, and the adjusted data as described above, on the right.

TABLE IV The S-factor of the ${}^2\text{H}(p,\gamma){}^3\text{He}$ at a few selected energies determined from data using Bayesian model averaging as described in the text. The first uncertainty is the statistical and the second is the model selection uncertainty. The value at any energy in the fitted range can be obtained by running the analysis provided at <https://github.com/nrp-g/leaner>.

$S_{12}(E)[10^{-7}\text{MeVb}]$								
$E = 0 \text{ keV}$	10	20	40	80	91	100	120	
2.028(51)(9)	2.644(60)(8)	3.276(70)(7)	4.579(94)(5)	7.31(15)(0)	8.11(16)(0)	8.77(18)(0)	10.24(21)(0)	

V. The ${}^3\text{He}({}^3\text{He}, 2p){}^4\text{He}$ reaction (S_{33})

The ${}^3\text{He}({}^3\text{He}, 2p){}^4\text{He}$ reaction terminates the pp-I chain. The ratio of its rate to that of the ${}^3\text{He}(\alpha, \gamma){}^7\text{Be}$ reaction controls the branching to the pp-II and pp-III chains, so that in SF I increasing S_{33} was discussed as a potential solution to the Solar Neutrino Problem. Subsequent experiments, notably a very low energy measurement at the LUNA 50 kV accelerator deep underground in Gran Sasso (Bonetti *et al.*, 1999) and a complementary experiment at somewhat higher energies (Kudomi *et al.*, 2004) ruled out such an increase in S_{33} , as summarized in SF II.

A. Shape of the particle spectrum

Since no new absolute measurements of S_{33} have been reported since SF II, we consider the same four experiments (Bonetti *et al.*, 1999; Junker *et al.*, 1998; Krauss *et al.*, 1987; Kudomi *et al.*, 2004). However, there is new information on the energy spectrum of the emitted protons. This spectrum has recently been measured using Inertial Confinement Fusion (ICF) plasmas for a Gamow peak energy of 165 keV (Zylstra *et al.*, 2017). The results show significant structure, indicating the presence of a sequential reaction mechanism passing through the unbound ground state of ${}^5\text{Li}$. This spectrum is impor-

tant for all of the S_{33} measurements since they determined cross sections by detecting only the protons above an energy threshold.

The efficiency correction by which Krauss *et al.* (1987) account for the threshold is not well documented, but the more recent measurements (Bonetti *et al.*, 1999; Junker *et al.*, 1998; Kudomi *et al.*, 2004) utilized the GENBOD event generator (James, 1968). It employs a simple reaction model without final state interactions, the Pauli principle, or Coulomb effects that are important near the spectrum endpoint. The only angular correlations are those required by energy and momentum conservation, and together these simplifications give simple ellipses for the singles energy distributions of the emitted nuclei. Published proton spectra obtained using accelerator beams do exist (Fig. 2 of Dwarakanath (1974) and Fig. 3 of Krauss *et al.* (1987)), and although they are not corrected for instrumental effects, they do not appear to be well described by ellipses.

The ${}^3\text{He}({}^3\text{He},2p){}^4\text{He}$ reaction may proceed via several sequential mechanisms, including $p + {}^5\text{Li}$ (with ${}^5\text{Li}$ in its ground or first excited state) and di-proton + α (where the di-proton is two correlated protons in a singlet state) (Brune *et al.*, 2015). More complicated three-body decay channels, sometimes called direct decays, are also possible. Only a relatively narrow intermediate state (here, the 1 MeV wide ${}^5\text{Li}$ ground state) could produce a peak in the energy spectrum, and precise classification of the reaction mechanism is in general both experimentally and theoretically ambiguous. When coincident detection of reaction products is used [as in Bonetti *et al.* (1999)], possible angular correlations between the reaction products also matter. A reaction through the $3/2^-$ ground state of ${}^5\text{Li}$ would emit the second proton preferentially either along or opposite the direction of the first proton, while di-proton emission would tend to send both protons in the same direction (Brune *et al.*, 2015).

The solid curves in Fig. 6 show two different R-matrix models of the proton energy spectrum at $E_{\text{c.m.}} = 165$ keV, fitted to the spectrum measured by Zylstra *et al.* (2017) for $E_p \gtrsim 6$ MeV. The measured spectrum differs significantly from the elliptical spectrum of GENBOD. Based on their Fig. 6, the experiment of Junker *et al.* (1998) had a detection threshold of about 5 MeV proton energy. The measurements by Bonetti *et al.* (1999) required a coincidence between two detectors, with a detection threshold of 2 MeV proton energy in each detector. Finally, Fig. 19 of Kudomi *et al.* (2004) indicates a detection threshold of about 4 MeV proton energy. We have estimated their sensitivities to the assumed proton spectrum by integrating the curves shown in Fig. 6 above energy thresholds of 2, 4, and 6 MeV. These integrals vary from 3% below to 6% above the result from an elliptical spectrum, depending on the specific threshold and the assumed spectrum.

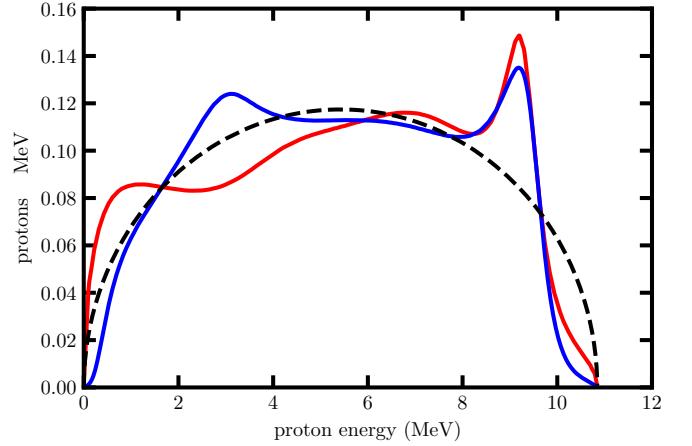


FIG. 6 Calculated proton energy spectra in the c.m. frame for the ${}^3\text{He}({}^3\text{He},2p){}^4\text{He}$ reaction at $E_{\text{c.m.}} = 165$ keV. The solid red and blue curves are models 1 and 2, respectively, from Fig. 3(a) of Zylstra *et al.* (2017). The black dashed curve is the elliptical spectrum. The spectra are normalized to unit area.

B. Recommendation

Based on these investigations, our recommended S_{33} fit is the one from SF II, but with an additional 4% systematic uncertainty typical of the 3% to 6% corrections to the extrapolated proton spectra estimated above to take the uncertain spectral shape into account. This gives

$$S_{33}^{\text{best}}(E) = 5.21 - 4.90 \left(\frac{E}{\text{MeV}} \right) + 11.21 \left(\frac{E}{\text{MeV}} \right)^2 \text{ MeV b}$$

$$\delta S_{33}(E) = \left[0.118 - 1.516 \left(\frac{E}{\text{MeV}} \right) + 14.037 \left(\frac{E}{\text{MeV}} \right)^2 - 15.504 \left(\frac{E}{\text{MeV}} \right)^3 + 71.640 \left(\frac{E}{\text{MeV}} \right)^4 \right]^{1/2} \text{ MeV b.}$$

Only the larger constant term in the uncertainty differs from the previous evaluation, and it is due to the spectral shape discussed above.

New measurements of this reaction could provide more accurate absolute cross sections by using lower energy thresholds and improve our understanding of the proton energy spectra and angular correlations. The analysis of new measurements should use Monte Carlo simulations considering a variety of plausible energy spectra and angular correlations, in order to estimate the sensitivity to these effects. We note finally that the proton energy spectrum may depend on $E_{\text{c.m.}}$, as this has been found to be the case for neutrons from the mirror reaction ${}^3\text{H}(t,2n)\alpha$ (Gatu Johnson *et al.*, 2018).

VI. The ${}^3\text{He}(\alpha, \gamma){}^7\text{Be}$ reaction (S_{34})

The ${}^3\text{He}(\alpha, \gamma){}^7\text{Be}$ reaction proceeds via non-resonant capture to the ground and 429 keV first excited states of ${}^7\text{Be}$. It has been studied experimentally using three main methods, with references given in subsections VI.A and VI.C: First, by detecting at least two of the three γ -rays from the reaction and taking the angular correlation with the alpha beam direction into account (prompt- γ method). Second, by detecting the induced ${}^7\text{Be}$ (Tilley *et al.*, 2002) radioactivity (activation method). Third, by counting the ${}^7\text{Be}$ recoils (recoil method).

A. Previous S_{34} recommendation in SF II

The ${}^3\text{He}(\alpha, \gamma){}^7\text{Be}$ S-factor recommended in SF II was developed in multiple steps. First, a model was selected for the shape of $S(E)$, based on existing nucleon-level calculations (Kajino, 1986; Nollett, 2001). A rescaling of the curve was fitted to data, but only at $E \leq 1.002$ MeV center-of-mass to minimize the role of short-distance physics in the models.

In SF I, a possible systematic discrepancy between data from the two methods previously used was discussed. By the time of SF II, in-depth studies from two groups using both activation and prompt methods were available, from LUNA (Bemmerer *et al.*, 2006; Confortola *et al.*, 2007; Gyürky *et al.*, 2007) and Seattle (Brown *et al.*, 2007) groups. These studies did not find any discrepancy between activation and prompt- γ data in direct comparisons. SF II opted to limit the fitting to the data by the activation and recoil methods. The prompt- γ data were left out because of their somewhat larger common mode errors, concerns about how well the γ -ray angular distribution was known, and avoidance of the correlated errors between activation and prompt data from the same experiment.

In SF II, only data published after 1998 were included. The SF II recommended value was $S_{34}(0) = 0.56 \pm 0.02_{\text{expt}} \pm 0.02_{\text{theor}}$, based on the data by the Weizmann (Singh *et al.*, 2004), LUNA (Bemmerer *et al.*, 2006; Confortola *et al.*, 2007; Gyürky *et al.*, 2007), Seattle (Brown *et al.*, 2007), and ERNA (di Leva *et al.*, 2009) groups.

B. Theory progress on S_{34}

Significant theoretical work on this reaction has occurred since SF II, but the basic understanding of its mechanism remains unchanged from the 1960s (Tombrello and Parker, 1963). It is dominated by external non-resonant capture into the two bound states, and most of the dipole strength at low energy arises beyond the range of nuclear interaction. The external

capture part of the cross section is determined by the ANC's of the bound states and by scattering phase shifts; near threshold, most of the strength lies at $\sim 5\text{--}20$ fm (Neff, 2011). In models with explicit wave functions the shorter-range strength largely cancels out due to effects of nucleon-exchange antisymmetry. All models feature a shallow minimum of $S(E)$ near 1.25 MeV, where capture from d -waves becomes comparable to that from s -waves.

The first fully *ab initio* calculation (Neff, 2011) appeared just after the SF II analysis concluded; it used the fermionic molecular dynamics (FMD) method and a softened representation of the Argonne v_{18} interaction (Wiringa *et al.*, 1995). It agrees well with both the scale and the energy dependence of the modern S-factor data, and it is very close to the energy dependence assumed in SF II. Neff's elastic-scattering phase shifts also agree well with experiment. Notably, Neff found that the dipole strength distribution departs significantly from pure external capture at ${}^3\text{He}$ - ${}^4\text{He}$ separations as large as 9 fm, compared with 3-4 fm in potential models.

Another *ab initio* model (Dohet-Eraly *et al.*, 2016; Vorabbi *et al.*, 2019) has also appeared, based on the no-core shell model with continuum (NCSMC) method and the chiral interaction of Entem and Machleidt (2003), “softened” by the similarity renormalization group (SRG) procedure. Freedom to choose the SRG stopping point allowed exact reproduction of the ${}^7\text{Be}$ breakup energy, which corrects for the main effect on the external capture of omitting the three-nucleon potential. The results agree well with the overall scale of the modern data, but their energy dependence and phase shifts depart from experiment, possibly due to the omission of explicit three-body forces.

An important advance since SF II has been the application of halo effective field theory (halo EFT) methods to astrophysical capture reactions (Bertulani *et al.*, 2002; Higa *et al.*, 2018; Premarathna and Rupak, 2020; Zhang *et al.*, 2020). EFTs are valid for systems with a natural separation between the momentum scales Q probed in low-energy experiment (e.g., corresponding to a binding energy or a projectile energy) and the much larger momentum scale Λ where the low-energy degrees of freedom are no longer valid: the word “halo” here refers to nuclei with only shallow bound states. In halo EFT the only degrees of freedom are the initial- and final-state nuclei treated as point particles, plus photons. For S_{34} , $\Lambda \sim 200$ MeV corresponds to the momentum needed to separate a proton from ${}^3\text{He}$ or to excite ${}^4\text{He}$; the thresholds to separate the two ${}^7\text{Be}$ bound states into ${}^3\text{He}$ and ${}^4\text{He}$ correspond to momenta $Q = 71.4$ MeV and 60.9 MeV.

Given a sufficient separation of scales, one constructs a Lagrangian that respects the system's symmetries and known qualitative features systematically, organized by powers of Q/Λ . This series can be truncated (at leading order or LO, next-to-leading order or NLO, next-to-next-to-leading-order or NNLO, etc.), and the precision

of the resulting theory depends on the sizes of the omitted terms. The precision of a calculation and the energy where it breaks down can be estimated by assuming coefficients of the first omitted term to have a “natural” size. Coupling constants of the Lagrangian must be fitted to data, and consistency of the power counting scheme (identification of powers of Q/Λ for the main operator terms of a system) is tested by whether the fitted constants have natural sizes. Low-order coupling constants in a halo EFT can often be identified with familiar quantities like ANC_s, scattering lengths, and effective ranges; the description of elastic scattering in halo EFT reproduces the (Coulomb-modified) effective-range expansion (Higa *et al.*, 2008; Kong and Ravndal, 2000; Poudel and Phillips, 2022). When a separation of scales exists, halo EFT is well-suited to data extrapolation because it avoids the tacit, hard-to-test, and unavoidable prior assumptions present in the short-range parts of models based on explicit wave functions.

Two groups have studied S_{34} in halo EFT (Higa *et al.*, 2018; Premarathna and Rupak, 2020; Zhang *et al.*, 2020). The sizes of the Coulomb interaction and the large s -wave scattering length make the correct power counting tricky to establish for this system. The work of Higa *et al.* (2018) and Premarathna and Rupak (2020) includes a careful examination of possible power-counting schemes and strict adherence to a power counting once established, up to terms of NLO or NNLO, respectively, in two different power-countings. These authors fitted EFTs to S-factor data, and they examined effects of including or excluding elastic scattering constraints in their fits. When they included scattering phase shifts from Boykin *et al.* (1972) in their fits, they concluded that the large-scattering-length power counting of the NLO theory was favored. This result promotes two-body currents (i.e., contributions not equivalent to external non-resonant capture) to leading order in the theory. The d -wave contribution that becomes important above 1 MeV first appears at NLO. In addition to data-fitting errors, these theories have errors estimated to be 10% from EFT truncation.

The work of Zhang *et al.* (2020) followed a different approach to power counting up through NLO. At $E < 2$ MeV, their derived expression for $S_{34}(E)$ is essentially the same as that from the Premarathna and Rupak (2020) NLO theory. Zhang *et al.* (2020) also developed some *ad hoc* (i.e., not systematically developed) higher-order EFT terms, referred to as partial-N4LO, to test for their impact on the fitting. The additional terms proved not to be required by the data and did not improve the fit; this result was taken to indicate that corrections from omitted terms are not large compared with experimental errors below 2 MeV. Scattering data were not considered in this work apart from very broad priors on scattering length and effective range; the correlated errors in the S-factor data were taken into account.

Like halo EFT, the phenomenological R-matrix approach avoids a model of nuclear interactions and uses a systematic parameterization to fit data (Descouvemont and Baye, 2010). At the time of SF II, S_{34} had been the subject of very little R-matrix fitting, apparently consisting only of the simple treatment in Descouvemont *et al.* (2004) that focused on BBN energies. After SF II an R-matrix analysis using the AZURE2 code (Azuma *et al.*, 2010; Uberseder and deBoer, 2015) was carried out on both elastic scattering and S-factor data in conjunction with the Notre Dame experiment (Kontos *et al.*, 2013), and a more elaborate analysis using Monte Carlo sampling to estimate errors was later reported in deBoer *et al.* (2014). In the latter analysis both the fitted value and error bars were heavily influenced by the numerous scattering data of Barnard *et al.* (1964). Very recently, the BRICK software package has been constructed to carry out Bayesian parameter estimation for AZURE2 and applied to both capture and elastic scattering at all energies in the ${}^7\text{Be}$ system by Odell *et al.* (2022b). In this work it was shown that markedly different $S_{34}(0)$ values result from inclusion or not of the older scattering data (Barnard *et al.*, 1964) alongside the very recent SONIK scattering data of Paneru *et al.* (2024).

A small amount of additional theoretical work in more traditional frameworks has appeared since SF II. This includes fits of potential models to S-factor data (Turtsunov *et al.*, 2018, 2021) and some calculations of the RGM type (Solovyev and Igashov, 2017, 2019).

C. Experimental progress on S_{34}

Since SF II, five new experiments have been reported: Four from the Madrid and ATOMKI groups by the activation technique, at relatively high center of mass energies (Bordeanu *et al.*, 2013; Carmona-Gallardo *et al.*, 2012; Szücs *et al.*, 2019b; Tóth *et al.*, 2023), and one from the Notre Dame group using the prompt- γ method (Kontos *et al.*, 2013).

Following the approach adopted in SF II, we again only use the recoil and activation data for the S_{34} fits below. The γ -ray angular distribution is not known experimentally (Brown *et al.*, 2007), and the resultant uncertainty increases the common-mode error for the prompt- γ studies somewhat. (An experiment to address previously-raised concerns about angular distribution (Brown *et al.*, 2007) has recently concluded at Felsenkeller Dresden. The data suggest a higher than expected anisotropy but are so far only available in the form of a PhD thesis (Turkat, 2023). If confirmed, they may lead to corrections of a few percent for some of the in-beam experiments where only one angle was instrumented.) In addition, all of the “modern” works reporting prompt- γ data except for Notre Dame (Kontos *et al.*, 2013) also include data obtained with other methods, leading to partial cor-

relations between data sets that would complicate fitting.

Following this restriction, our data selection proceeds as follows. First, the four data sets previously used in SF II are carried over here: Weizmann (Singh *et al.*, 2004), LUNA (Bemmerer *et al.*, 2006; Confortola *et al.*, 2007; Gyürky *et al.*, 2007) (only the activation data), Seattle (Brown *et al.*, 2007) (only the activation data), and ERNA (di Leva *et al.*, 2009).

Two of the archival data sets excluded by the 1998 cutoff date in SF II merit some further discussion here: The data of Osborne *et al.* (1982, 1984) consist of two points measured by activation in a ${}^3\text{He}$ gas cell. However, their uncertainties are not separated into statistical and systematic components. Including them would have required an uncertain guess on how to divide the error bars, and the fit result would barely change, so we left them out. The work by Robertson *et al.* (1983) consists of one data point that was obtained by averaging two separate activation measurements with a ${}^3\text{He}$ and a ${}^4\text{He}$ gas cell, respectively. We do not use this result because we lack details to verify the background subtraction, an issue that was raised during our meeting by one of its authors. We also choose not to use the activation study by Volk *et al.* (1983). In that experiment there were thick entrance foils and the ${}^4\text{He}$ beam was completely stopped inside a high-pressure ${}^3\text{He}$ gas cell. This gave an integrated measurement over a wide energy range, so that the analysis depends strongly on the assumed shape of the S_{34} curve, and the result was reported only as an extrapolated $S_{34}(0)$.

We now consider the new data since SF II. An activation experiment in Madrid (Carmona-Gallardo *et al.*, 2012) used a ${}^3\text{He}$ beam incident on a ${}^4\text{He}$ gas cell, reporting three data points. Another activation experiment was reported by the ATOMKI group, using a ${}^4\text{He}$ beam on ${}^3\text{He}$ gas cells (Bordeanu *et al.*, 2013). Two higher-energy campaigns at ATOMKI were again performed using the activation method. These latter data are at $E = 2.5\text{-}4.4$ MeV (Szücs *et al.*, 2019b) and $E = 4.3\text{-}8.3$ MeV (Tóth *et al.*, 2023), respectively, above the energy range suitable for halo EFT, and therefore not included.

A detailed study by the Notre Dame group reported 17 data points (Kontos *et al.*, 2013), using the primary γ -ray from ground state capture and the secondary γ -ray from the deexcitation of the 429 keV first excited state of ${}^7\text{Be}$. No activation data are reported in that work, and the γ -ray detector was placed at just one angle, 90° . Since only prompt- γ data are reported, we did not include this data set. In order to test the effects of this decision, we repeated some of our fits (see below, Section VI.D) with the modern prompt- γ experiments included: the 17 points from Kontos *et al.* (2013), 3 prompt- γ points from LUNA (Confortola *et al.*, 2007) and 8 from Seattle (Brown *et al.*, 2007). The extrapolated S-factor changed by less than 1%, well within the error bars for the recommended S_{34} .

Finally, an indirect experiment using the ${}^6\text{Li}({}^3\text{He},\text{d}){}^7\text{Be}$ reaction and the ANC technique has recently been reported (Kiss *et al.*, 2020). It is left out of our fits due to the additional normalization and theory uncertainties involved in determining an ANC from a transfer experiment, which are larger than the errors in the S-factor measurements used for the present fits. All the data used in the present fits are summarized in Table V.

As in SF II, we model data uncertainties as consisting of a component that is independent for each point and a common-mode component that applies to all data from a given experiment as a multiplicative factor. This separation is well-documented for all of the modern data and is also shown in Table V. Except at the lowest energies, the common-mode error typically dominates.

D. Data fitting

We base our recommended S_{34} on fits to the halo EFT and R-matrix parameterizations discussed above. These avoid tacit assumptions present in potential-models and conceptual difficulties involved in combining *ab initio* constraints with data. The fits presented here differ from the previously published fits of Premarathna and Rupak (2020), Zhang *et al.* (2020), and Odell *et al.* (2022b) mainly in the uniform use of the agreed-upon capture data and uncertainties from Section VI.C across all fits. We restricted fitting to the $E < 2$ MeV range of validity for the NLO halo EFT expressions. Despite uniform handling of capture data, the fits in each framework handle scattering inputs differently for reasons discussed below.

In addition to the total capture cross section and scattering data, the fitted data also include branching ratios for capture into the two ${}^7\text{Be}$ bound states, taken from Brown *et al.* (2007), di Leva *et al.* (2009), Confortola *et al.* (2007), and Kontos *et al.* (2013). These are necessarily from prompt- γ experiments and suffer from the concerns about angular distribution discussed in Section VI.C. However, their inclusion simplifies the fitting considerably by breaking parameter degeneracies between ground- and excited-state transitions (especially in the fitted ANCs), probably without strong impact on $S_{34}(0)$.

In constructing fits we split common-mode and point-to-point errors and “float the norms” of data sets using the cost function

$$\chi^2 = \sum_{\alpha=1}^{N_{\text{set}}} \sum_{i=1}^N \frac{(y_i - \mu_i/s_{\alpha})^2}{\sigma_{\alpha,i}^2} + \sum_{\alpha=1}^{N_{\text{set}}} \frac{(1 - s_{\alpha})^2}{\omega_{\alpha}^2} \quad (13)$$

to describe goodness of fit (D’Agostini, 1994).

A rescaling factor s_{α} is fitted for each data set α , with its deviation from unity penalized by the common-mode errors ω_{α} given in Table V. The index i sums over all

TABLE V Experimental data used for the S_{34} fit. See text for details. For each experimental data set, the rescaling factors s_α are determined for the Higa *et al.* (2018), Zhang *et al.* (2020), and BRICK ((Odell *et al.*, 2022b)) without or with the inclusion of elastic scattering.

Group and references	Energy range [keV]	Data points			Common mode		Rescaling factor s_α		
		used	total	uncert.	Higa EFT	Zhang EFT	BRICK	BRICK	BRICK+S
Weizmann (Singh <i>et al.</i> , 2004)	420–950	4	4	2.2%	1.03(2)	1.02(2)	1.02(2)	1.03(2)	
LUNA (Bemmerer <i>et al.</i> , 2006; Confortola <i>et al.</i> , 2007; Gyürky <i>et al.</i> , 2007)	93–170	7	7	3.0%	1.02(2)	1.04(2)	1.05(2)	1.01(2)	
Seattle (Brown <i>et al.</i> , 2007)	327–1235	8	8	3.0%	0.96(1)	0.95(2)	0.95(2)	0.98(2)	
ERNA (di Leva <i>et al.</i> , 2009)	650–2504	27	27	5.%	0.96(3)	0.94(2)	0.96(2)	0.99(2)	
Madrid (Carmona-Gallardo <i>et al.</i> , 2012)	1054–2804	2	3	5.2%	0.99(3)	0.99(3)	0.97(2)	0.99(2)	
ATOMKI (Bordeanu <i>et al.</i> , 2013)	1473–2527	4	5	5.9%	1.01(3)	1.00(3)	1.01(3)	1.04(3)	

points within a given data set; y_i is a measured cross section, μ_i is a predicted cross section, and $\sigma_{\alpha,i}$ is the point-to-point error of the i th point in data set α .

We carried out both frequentist fits that minimize χ^2 and Bayesian fits based on the posterior probability distribution of a likelihood function computed from $\exp(-\chi^2/2)$. In the Bayesian fits parameters, extrapolated $S_{34}(0)$, and their errors were determined by Monte Carlo sampling of the parameter space. The relatively large number of model parameters and the significant fitting degeneracies between some of them make Bayesian analysis a natural choice for finding best fit and confidence intervals for the multi-parameter models applied here. Two of the groups ran into serious difficulty with the frequentist fits due to parameter degeneracies and shallow local minima; during this study they were unable to produce frequentist fits in which they had confidence.

We performed multiple fits in the NLO halo EFT of Higa *et al.* (2018) and Premarathna and Rupak (2020). These fits were carried out both with and without scattering constraints, which mainly impact $S_{34}(0)$ by removing parameter degeneracies that would otherwise leave the s -wave scattering length poorly constrained. The code base for this version of halo EFT incorporates scattering data through phase shifts. These were taken from the partial-wave analysis of Boykin *et al.* (1972), which at low energy are based mainly on the data of Barnard *et al.* (1964).

One set of fits for the Higa *et al.* (2018) halo EFT was produced by χ^2 minimization, proceeding in two steps: first LO parameters were fitted to $E \leq 1000$ keV data, and then the results were taken as initial values in the search to minimize χ^2 over all parameters in the NLO theory for the full set of data. These fits gave $S_{34}(0) = 0.566 \pm 0.025$ keV b with the scattering constraint and $S_{34}(0) = 0.588 \pm 0.015$ keV b without (uncertainties being propagated in linear approximation using covariances and partial derivatives). Formally the EFT truncation error from stopping at NLO corresponds to an additional theoretical error that can shift S-factors by 10%. However, any fitted curve is constrained by low-energy data, which in some sense become effectively renormalization conditions of the field theory; the error

on extrapolated $S_{34}(0)$ should probably be smaller than 10% by an amount that is hard to estimate.

A second set of fits to the Higa *et al.* (2018) halo EFT was carried out using Bayesian methods. Priors for the EFT parameters were developed based on previous experience, and data rescaling factors were incorporated as additional priors. The Bayesian results (including only experimental error) are $S_{34}(0) = 0.561^{+0.017}_{-0.018}$ keV b including the phase shifts by Boykin *et al.* (1972) and $S_{34}(0) = 0.559^{+0.018}_{-0.019}$ keV b excluding them, consistent with the χ^2 -minimization.

Searches of the parameter space to minimize χ^2 for the NLO halo EFT of Zhang *et al.* (2020) ran into difficulties with parameter degeneracy and local minima in the χ^2 surface; for this formalism we report only Bayesian results. No experimental information about scattering was used, but flat priors on scattering length and effective range were chosen over a 5σ range around recent experiment. This fit differs from Zhang *et al.* (2020) mainly by excluding the Notre Dame data and including the ERNA activation data, and it gives $S_{34}(0) = 0.581 \pm 0.016$ keV b. The error associated with EFT truncation at NLO is estimated in this approach by separately fitting the partial-N4LO theory discussed above. The result suggests that EFT truncation at NLO affects extrapolation from the data to threshold by ~ 2 –3%.

Our R-matrix fits are based mainly on sampling Bayesian posterior probabilities with the BRICK code. We also produced frequentist fits, but we were unable to estimate their errors convincingly. The R-matrix fits used capture data both alone and in combination with the SONIK scattering data (Paneru *et al.*, 2024) – the latter being chosen because of concerns with the Barnard *et al.* (1964) data that are discussed in Odell *et al.* (2022b). Since BRICK fits elastic differential cross sections directly, it was not feasible to use the same phase-shift-based scattering constraints as our halo EFT fits. The R-matrix fit that includes scattering data is essentially the “CS” fit of Odell *et al.* (2022b), but with a restriction to only the capture data described in Table V, only scattering data below 2 MeV center-of-mass, and only the R-matrix parameters relevant below 2 MeV (no 7/2–

TABLE VI Main results of the Bayesian fits performed on the S_{34} data. The ANC refers to the sum of the ANCs for the ground and excited state transitions. See text for details.

Fit	$S_{34}(23)$ [keV b]	S'_{34}/S_{34} [MeV $^{-1}$]	a_0 [fm]	C^2 [fm $^{-1}$]
Higa <i>et al.</i> (2018) EFT-NLO	0.554	-0.56	30^{+5}_{-2}	$26.9^{+4.5}_{-5.1}$
Zhang <i>et al.</i> (2020) EFT-NLO	0.573	-0.61	$46(6)$	29(3)
BRICK R-matrix	0.562	-0.57		$24.4(9)$
BRICK R-matrix+scattering	0.531	-0.47		$22.6(7)$
ANC (Kiss <i>et al.</i> , 2020)				34 ± 1.6

level or radiative widths for d -wave background poles).

Several of the fits are compared in Fig. 7. There it is apparent that they all agree within $\leq 5\%$ in the 0–2000 keV energy range, and within $\leq 4\%$ in the astrophysically relevant range 0–500 keV. The fit to the Higa *et al.* (2018) NLO EFT lies in the middle of the range, so we adopt it as a reference in the lower panel of the figure. Relative to that fit, the fits to the NLO EFT of Zhang *et al.* (2020) and the R-matrix fit without scattering rise up as much as 4% higher at 0–200 keV.

The most striking difference among the curves is between the BRICK R-matrix fit including the SONIK scattering data (Paneru *et al.*, 2024) and all the other fits. While the R-matrix fit with scattering finds rescaling factors somewhat closer to unity than the other fits do, its energy dependence in the 0–1000 keV range is qualitatively different, as is especially visible in the lower panel of Fig. 7 where residuals relative to the fitted Higa *et al.* (2018) EFT are shown. The BRICK fit without the scattering data is much closer to all the halo EFT fits, so the recent scattering data apparently have a large impact.

E. Recommended S_{34} value

Because the fit to the NLO EFT of Higa *et al.* (2018) gives the central result among those attempted here, we adopt it for our recommended $S(E)$. For ease of adoption, it is noted that this recommended S_{34} curve can be empirically parameterized using the same shape as in SF II, by the numerical equation (following a customary form in the past literature (Adelberger *et al.*, 2011; Kajino *et al.*, 1987))

$$S_{34}(E) = (0.5610 \text{ keV b}) \exp(-0.5374 E) \quad (14) \\ \times [1 - 0.4829 E^2 + 0.6310 E^3 - 0.1527 E^4].$$

Here E is the center of mass energy in MeV. Equation (14) reproduces our recommended curve within 0.3% and is only applicable for $E = 0$ –1600 keV.

Propagation of uncertainties from measurement to extrapolated $S_{34}(0)$ is relatively straightforward and unambiguous (at least in the Bayesian fitting). However, the uncertainty due to theoretical approximation is more

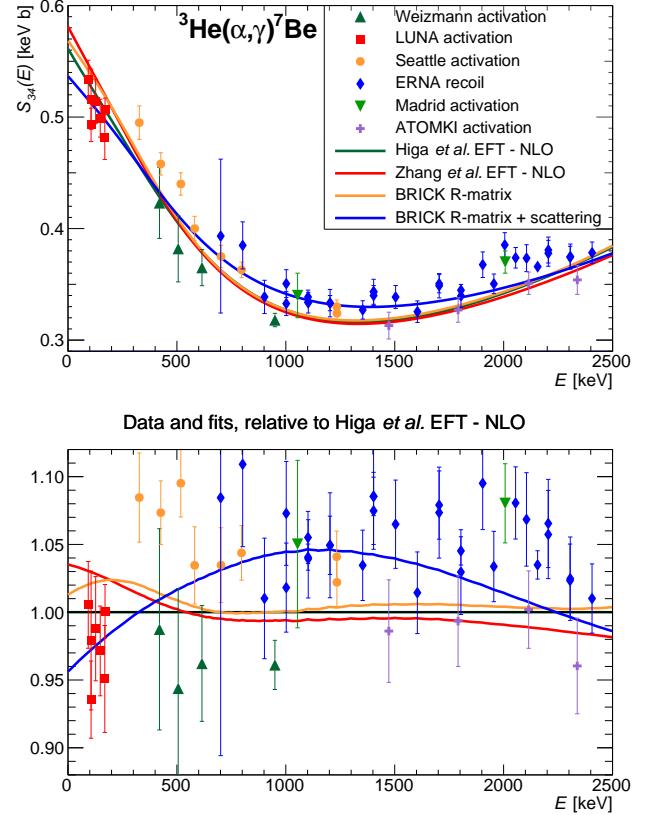


FIG. 7 Top panel: ${}^3\text{He}(\alpha, \gamma){}^7\text{Be}$ astrophysical S-factors from the experiments used for the fitting of the excitation function: Weizmann (Singh *et al.*, 2004), LUNA (Bemmerer *et al.*, 2006; Confortola *et al.*, 2007; Gyürky *et al.*, 2007), Seattle (Brown *et al.*, 2007), Madrid (Carmona-Gallardo *et al.*, 2012), and ATOMKI (Bordeanu *et al.*, 2013). The fit curves from two different EFT based fits (approaches based on Higa *et al.* (2018) and Zhang *et al.* (2020)) and from two R-matrix fits (BRICK Odell *et al.* (2022b), without or with elastic scattering data) are also shown. Bottom panel: Experimental data and fits, normalized to the Higa *et al.* (2018) based EFT fit. See Table V and text for details.

complicated, and error estimation on extrapolated quantities in EFTs remains an open area of research; estimates from the groups working on S_{34} are discussed above. We estimate the overall theoretical uncertainty on our recommended S-factor to be the measured by the 4% dispersion among fitted curves; this is comparable to the more formal estimates.

Our recommendations for the zero energy astrophysical S-factor and the low-energy slope are then

$$S_{34}(0) = 0.561 \pm 0.018_{\text{exp}} \pm 0.022_{\text{theor}} \text{ keV b} \quad (15)$$

$$\frac{S'_{34}(0)}{S_{34}(0)} = -0.54 \pm 0.07 \text{ MeV}^{-1}. \quad (16)$$

The given value of S'_{34}/S_{34} applies over $E = 0$ –500 keV. Its considerable uncertainty is given by the full span of slopes among the four fits discussed above, which is dom-

inated by the inclusion of elastic scattering in one of the R-matrix fits. The $S_{34}(E)$ curve is quite straight at low energy in all models; its second derivative is consistent with zero and even difficult to compute for a given fit because of numerical cancellations.

The first error bar for $S_{34}(0)$ is given by the fit uncertainty, conservatively using the largest value among the various fits, and it encodes the experimental uncertainty. The second error bar reflects the 4% dispersion among theoretical formalisms. The two errors can be added in quadrature for applications. Our recommendation is essentially the same as in SF II, but derived from completely different theoretical approaches that are much more directly connected to the data and much less dependent on model assumptions. It is also based on data over a wider energy range. The dispersion of S_{34} and S'_{34} among models at the solar Gamow peak is illustrated in Table VI, along with some indication of how fitted scattering lengths a_0 and squared ANCs C^2 vary among frameworks. (The latter are highly degenerate with short-range parameters in the fitting; in the Higa *et al.* (2018) EFT they are sensitive to small variations in the p -wave effective ranges that are fitted instead of ANCs.)

F. Recommendations for future work on S_{34}

Here we present recommendations for further work that would improve understanding of the ${}^3\text{He}(\alpha, \gamma){}^7\text{Be}$ reaction. Our first recommendation for experiment concerns elastic ${}^3\text{He} + {}^4\text{He}$ scattering. New experiments over a wide energy range including at least part of the astrophysical range are needed to address the marked impact of the recent SONIK scattering data (Paneru *et al.*, 2024) on fitting, for independent confirmation. Data below 2 MeV would be most useful for consistent fitting of halo EFTs.

Second, new measurements (besides the preliminary data of Turkat (2023)) are needed for the angular distribution of the emitted γ -rays. This would relieve concerns about including prompt- γ data, most notably the Notre Dame (Kontos *et al.*, 2013) that do not include activation measurements, in future fits. These experiments should also be designed to provide new angle-corrected branching ratios for capture into the two bound states.

Our third recommendation is a study of the astrophysical S-factor over a wide energy range, 50-1500 keV and extending to even lower energies when technically feasible, with small point-to-point errors. Even with a large common-mode error, this would significantly constrain theoretical curve shapes near threshold, and providing an independent check on the influence of scattering data on $S(E)$.

On the theoretical side, further development of *ab initio* reaction calculations is desirable, improving on points

of current difficulty like explicit inclusion of three-body potentials. Calculations with the same potential but by different methods would help to separate computational issues from the impact of potential choice. Further exploration of how *ab initio* inputs like ANCs and scattering lengths could serve as fitting priors would also be a useful avenue of research.

Authors of future theoretical calculations should consider including γ -ray angular distributions in their results; until very recently angular distributions were only available from two potential models (Kim *et al.*, 1981; Tombrello and Parker, 1963). New calculations would inform analysis of experiments, and the angular distribution measurements we recommend would provide tests of the calculations (or fitting constraints).

A source of complication here has been the need for a good meeting point between scattering data and theory. Given the difficulties of constructing a unique phase shift analysis without an underlying theory, and the recent arrival of modern scattering data, theorists should consider including elastic differential cross sections directly as outputs and fitting constraints.

VII. The hep reaction (S_{hep})

The reaction $p + {}^3\text{He} \rightarrow {}^4\text{He} + e^+ + \nu_e$, also known as the hep reaction, is one of the possible processes taking place after the ${}^2\text{H}(p, \gamma){}^3\text{He}$, and it produces the most energetic neutrinos, with an endpoint energy of 18.8 MeV. It is however the least probable, with its S-factor, and consequently its rate, even beyond the reach of current experiments. This is due to the fact that the hep reaction is induced by the weak interaction and further suppressed by Coulomb barrier and by the fact that the leading one-body axial (Gamow-Teller) current operator cannot connect the main s -state components of the $p + {}^3\text{He}$ and ${}^4\text{He}$ initial- and final-state wave functions. As a consequence, the one-body axial current non-zero contribution is due to the small components of the ${}^3\text{He}$ and ${}^4\text{He}$ wave functions. Furthermore, other contributions, such as weak magnetism and other one-body corrections to the vector current, and two-body vector and axial currents, and the p -waves in the $p + {}^3\text{He}$ initial state, normally suppressed at solar temperatures, are significantly enhanced. This is further complicated by cancellations between the one-body and two-body axial current contributions. As a consequence, similarly to the pp fusion, the hep reaction cross section is too small to be measured in laboratory and only theoretical predictions are available.

The most recent studies of this reaction, already evaluated in SF II, are those of Marcucci *et al.* (2000) and Park *et al.* (2003). The first one was performed within a phenomenological approach, similar to the calculation of Schiavilla *et al.* (1998) for the pp fusion. In this particular case, the $p + {}^3\text{He}$ and ${}^4\text{He}$ initial and final nuclear

wave functions were obtained with the correlated Hyper-spherical Harmonics (CHH) variational method (Viviani *et al.*, 1995, 1998), using the Argonne v_{18} (AV18) two-nucleon potential (Wiringa *et al.*, 1995) augmented by the Urbana IX (UIX) three-nucleon interaction (Pudliner *et al.*, 1995). In this way, the binding energies of ${}^3\text{He}$ and ${}^4\text{He}$, and the singlet and triplet $p+{}^3\text{He}$ scattering lengths were in good agreement with experiment. The weak vector and axial transition operators were obtained within a phenomenological approach. In particular, the largest two-body contribution to the axial current arising from the excitation of intermediate Δ -isobar degrees of freedom was included using the transition-correlation operator scheme of Carlson *et al.* (1991); Schiavilla *et al.* (1992) and fixing the nucleon-to- Δ axial coupling constant to reproduce Γ_β^T as in the pp case (Schiavilla *et al.*, 1998).

The second study of Park *et al.* (2003) was performed within the same hybrid EFT approach used for the pp fusion (EFT*), with the same nuclear weak current operators obtained within EFT as discussed in Section III, the LEC \hat{d}^R being fixed to reproduce Γ_β^T . The initial and final wave functions were derived using the CHH method with the AV18/UIX potential, as in Marcucci *et al.* (2000).

Following the results of Marcucci *et al.* (2000) and Park *et al.* (2003), SF II recommended for the zero-energy S-factor, $S_{\text{hep}}(0)$, the value $(8.6 \pm 2.6) \times 10^{-20}$ keV b. Three recommendations were also made: (i) to perform new studies with a broad spectrum of Hamiltonian models, in order to properly access the theoretical uncertainty; (ii) to study other weak reactions for which experiments can test the predictions made by employing the same theoretical ingredients as those used for the hep reaction calculation; (iii) to further understand the relation between the hep reaction and the so-called hen process ($n + {}^3\text{He} \rightarrow {}^4\text{He} + \gamma$). Of these recommended works, only the second one has been performed so far, applying the theoretical framework used for the hep study to the muon capture reactions on light nuclei (Marcucci *et al.*, 2011).

Given the lack of further studies on the hep reaction after SF II, we decide to maintain

$$S_{\text{hep}}(0) = (8.6 \pm 2.6) \times 10^{-20} \text{ keV b} \quad (17)$$

as recommended value. Notice that we could have increased $S_{\text{hep}}(0)$ by $\sim 1\%$, in analogy with $S_{11}(0)$, due to the updated values of input parameters (mostly G_A/G_V). However, we have decided not to apply such an increase, because it does not derive from a direct calculation, and, most of all, because such an increase lies well within the quoted $\sim 30\%$ uncertainty. However, given the advances in *ab initio* studies of the last decade and in the development of the EFT framework, we update the recommendations of SF II as follows: (i) the hep process should be

revisited within a fully consistent χ EFT approach, similarly to what has been done in these years for the pp fusion and other processes, as, for instance, muon captures on light nuclei (Acharya *et al.*, 2018; Bonilla *et al.*, 2023; Ceccarelli *et al.*, 2023; Gnech *et al.*, 2024; Marcucci *et al.*, 2012, 2018); (ii) the initial and final wave functions should be calculated with the most accurate *ab initio* methods, as the (uncorrelated) HH method (Kievsky *et al.*, 2008; Marcucci *et al.*, 2020), which has been proven to provide more accurate $A = 4$ bound and scattering wave functions than the CHH method, using both local and non-local interactions; (iii) the relation between the hep and the hen process should be understood. The hen process has been studied within χ EFT and the HH method by Viviani *et al.* (2022), but a consistent parallel study of both $A = 4$ reactions is still missing.

A relatively recent SNO Collaboration analysis of data from all three of the detector's running phases yielded a one-sided confidence-interval bound on the hep neutrino flux of Aharmim *et al.* (2020)

$$\phi_{\text{hep}} < 30 \times 10^3 \text{ cm}^{-2}\text{s}^{-1} \text{ (90\% C.I.)}, \quad (18)$$

a result in agreement with the SSM prediction. The lack of a definite measurement, however, has meant that there is no substantive experimental test of the predicted $S_{\text{hep}}(0)$. Various global analyses of solar neutrino data have provided weak evidence for a nonzero flux, e.g., from Bergstrom *et al.* (2016)

$$\phi_{\text{hep}} \Big|_{\text{global analysis}} = 19^{+12}_{-9} \times 10^3 \text{ cm}^{-2}\text{s}^{-1}, \quad (19)$$

which is consistent with the experiment bound given above.

An extraction of the hep flux from experiment will likely require a detailed shape analysis to separate hep neutrinos from the high-energy tail of the ${}^8\text{B}$ spectrum and the low-energy tail of atmospheric neutrinos. Typically the hep spectrum is assumed to have an allowed shape: indeed, in the work reported here spectra are not provided, where p -waves, weak magnetism, and similar corrections have been included. Because these corrections are unusually large for hep neutrinos, future studies should provide spectra and evaluate the impact of such corrections on the shape.

VIII. Electron capture by $p+p$ and ${}^7\text{Be}$

Electron capture reactions are the sources of lines found in the solar neutrino spectrum. In particular, electron capture on $p + p$, i.e., the process $p + p + e^- \rightarrow d + \nu_e$, also known as the pep reaction, competes with $p + p$ fusion and depends on the same nuclear matrix element. Therefore, the ratio between the pp and pep rates

is independent of nuclear physics. Based on this consideration, in **SF II** the result of **SF I**, which is based on the work of Bahcall and May (1969), was multiplied by the radiative corrections calculated by Kurylov *et al.* (2003), leading to the final result given by

$$R(\text{pep}) = 1.130(1 \pm 0.01) \times 10^{-4} (\rho/\mu_e) \times T_6^{-1/2} [1 + 0.02(T_6 - 16)] R(pp), \quad (20)$$

where ρ is the density in units of g/cm³, μ_e is the mean molecular weight per free electron, and T_6 is the temperature in units of MK. The range of validity is $10 < T_6 < 16$. Given that no new evaluation of the pep rate has been performed since **SF II**, the result of Eq. (20) represents also the present recommended value.

Competition between electron and proton capture on ⁷Be fixes the branching ratio of the ppII and ppIII chains, and thereby the ⁷Be and ⁸B neutrino fluxes. Supralow-allowed electron capture on the $J^\pi = 3/2^-$ ground state

of ⁷Be leads to the $J^\pi = 3/2^-$ ground state or the $J^\pi = 1/2^-$, 478 keV first excited state of ⁷Li with a measured (terrestrial) branching ratio of 10.44(4)% (Tilley *et al.*, 2002). It is customary to calculate the solar decay rate in terms of the measured terrestrial decay rate, which has the advantage of removing the nuclear physics dependence, but it requires a proper calculation of the densities of electrons at the nucleus. This has usually been done by considering continuum and bound electrons separately, using the Debye-Hückel (DH) approximation pioneered by Salpeter (1954) for the screening of bound electrons in the solar plasma; the capture rate of continuum electrons is not significantly influenced by electron screening. Furthermore, it is necessary to calculate the atomic probability densities governing the K and L terrestrial electron capture rates. Fretwell *et al.* (2020) recently measured the ratio of L to K capture using superconducting tunnel junctions. The recommendation of both **SF I** and **SF II** was

$$R(^7\text{Be} + e^-) = 5.60(1 \pm 0.02) \times 10^{-9} (\rho/\mu_e) T_6^{-1/2} [1 + 0.004(T_6 - 16)] \text{ s}^{-1}, \quad (21)$$

which is valid for $10 < T_6 < 16$. This expression is based on the continuum capture rate calculated by Bahcall and Moeller (1969), the average ratio of the total capture rate to the continuum rate calculated using three solar models in Bahcall (1994), and the current value of the half-life, 53.22(6) d (Tilley *et al.*, 2002). It agrees within 1% with a density matrix calculation of the rate that makes no assumptions regarding the nature of electronic quantum states in the solar plasma and allows for aspherical fluctuations in the spatial distributions of plasma ions (Gruzinov and Bahcall, 1997). The estimated uncertainty of 2% accounts for possible corrections to the Debye-Hückel approximation due to thermal fluctuations in the small number of ions in the Debye sphere, and breakdowns in the adiabatic approximation (Johnson *et al.*, 1992).

Although calculations have appeared in the literature claiming new plasma effects based on model calculations (see Simonucci *et al.* (2013); Vescovi *et al.* (2019)), we continue to regard the arguments in Bahcall *et al.* (2002) as definitive. That reference presented five distinct derivations demonstrating that the Salpeter formula for screening corrections is valid for pp-chain reactions and solar conditions, up to corrections on the order of a few percent.

IX. The ${}^7\text{Be}(p, \gamma){}^8\text{B}$ reaction (S_{17})

A. Introduction

Radiative proton capture on ⁷Be ($J^\pi=3/2^-$) at solar energies proceeds through non-resonant capture to the ground state of ⁸B ($J^\pi=2^+$). This capture occurs predominantly at separations well beyond the range of the strong interaction via $E1$ transitions from s and d partial waves, as the $M1$ amplitude and contributions from higher partial waves are negligible in the energy range of interest (Christy and Duck, 1961). The importance of this reaction for the determination of the high energy solar neutrino spectrum and the experimental data reviewed by **SF I** inspired several experiments that, taken together, provided a consistent picture of the energy dependence of the reaction cross section, despite discrepancies in absolute scale. Prior to the review of **SF II**, most direct measurements were performed using a radioactive target and an intense proton beam. In spite of the great care taken in evaluating the systematic energy-independent uncertainties, in the following referred to as common-mode errors (CMEs), a remarkable discrepancy persisted even in the most recent experiments, that was finally handled in **SF II** by inflating the stated experimental uncertainties. Indirect measurements of the cross section of ${}^7\text{Be}(p, \gamma){}^8\text{B}$ based on Coulomb dissociation were considered, but not included in the final recommendation of **SF II** for the zero-energy astrophysical S-factor $S_{17}(0)$. However, despite the experimental issues, the uncertainty

on the recommended value of $S_{17}(0)$ in **SF II** is dominated by the theoretical contribution. Therefore, our analysis focused on these aspects (discrepancy of experimental data, indirect measurements, theoretical uncertainties), as discussed in the next sections.

B. Experimental Data

1. Direct Measurements

The thorough review of published work done in **SF II** investigated the influence of beam-target overlap, target stoichiometry, beam energy loss, and the backscattering of ${}^8\text{B}$ recoils on CMEs and led to the selection of a homogeneous group of well-documented data sets. Since then, no further information became available on these aspects of direct measurements. The data that formed the basis of the **SF II** recommendation were those of Baby *et al.* (2003a), Baby *et al.* (2003b), Hammache *et al.* (1998), Hammache *et al.* (2001), Hass (1999), Strieder *et al.* (2001), and the BE3 data set of Junghans *et al.* (2003). We considered the complete set of radiative capture measurements that were the basis of the **SF II** recommendation for $S_{17}(0)$, where we adopted the revised BE3 data presented by Junghans *et al.* (2010), which were reanalyzed using a more sophisticated model of the target and more accurate stopping powers. In addition, we included in our analysis the only new experiment reported thereafter of Buompane *et al.* (2022), which used a radioactive ${}^7\text{Be}$ ion beam and the recoil mass separator ERNA to detect the ${}^8\text{B}$ recoils. While these new results contribute to the determination of S_{17} with different, well controlled systematics (Buompane *et al.*, 2018), their large statistical uncertainty relative to the other experiments, due to the low counting statistics, limits their impact.

New scattering data have been published. Elastic and inelastic scattering cross sections at relative kinetic energies between 474 keV and 2.74 MeV were measured and s wave scattering lengths were inferred by Paneru *et al.* (2019). At higher energies between 1.6 and 3.4 MeV, thick target elastic and inelastic scattering excitation function measurements were used to infer the existence of new resonances above 1.8 MeV (Mitchell *et al.*, 2013). These high lying resonances have no direct influence on $S_{17}(0)$ because of their small partial widths in the ${}^7\text{Be}_{\text{gs}} + p$ channel and/or because they cannot be populated by s - or d -wave capture at low energy.

2. Indirect Measurements

In **SF II**, Coulomb dissociation measurements based on the formalism developed by Baur *et al.* (1986) were considered (Davids *et al.*, 2001a,b; Iwasa *et al.*, 1999; Kikuchi *et al.*, 1998; Kikuchi *et al.*, 1997; Motobayashi *et al.*, 1994; Schümann *et al.*, 2003; Schümann *et al.*, 2006), where in

some cases the most recent publications supersede the previously published ones. Finally, the results of these experiments were not included in the determination of $S_{17}(0)$ due to disagreements over whether the analyses of the various measurements properly accounted for the contributions of $E2$ transitions. We disagree with these reservations and find the agreement of these measurements with radiative capture measurements significant. However, in order to include these consistent data in formulating our recommendation for $S_{17}(0)$, we inflated the CMEs of the Coulomb breakup measurements to adequately account for the uncertainty in the $E2$ components assumed in each analysis.

The common mode errors are obtained from a linear sum of those given in the manuscripts and an additional common mode error due to the estimated size of the $E2$ component in each measurement. This additional error is applicable in one direction only, downward for the Kikuchi *et al.* (1998) and Schümann *et al.* (2006) analyses of the RIKEN and GSI measurements, respectively, and upward for the Davids and Typel (2003) analysis of the Michigan State University (MSU) measurement, according to whether the $E2$ component was included or excluded. Though the Davids and Typel (2003) analysis did include a common mode error contribution of 2.5% due to the $E2$ component (which represented 5% of the total cross section), we have included an additional 2.5% contribution to the MSU experiment. The size of the additional common mode errors for the RIKEN and GSI measurements was based on scaling this 2.5% according to the predicted size of the $E2$ contribution to the cross section in these measurements relative to that in the MSU measurement, evaluated using first-order perturbation theory and the $E1$ and $E2$ S-factors calculated in a potential model of ${}^8\text{B}$ (Bertulani, 1996). Thus the additional single-sided common mode errors added to the GSI, MSU, and RIKEN measurements are -2.1%, 2.5%, and -4.3%, respectively.

C. Theory

The size of the theoretical contribution to the uncertainty in the **SF II** estimate of $S_{17}(0)$ was twice that of the experimental contribution. Since then, significant progress has been made. A new potential model calculation was performed by Dubovichenko *et al.* (2019). The large-scale computational demands of calculations using high-quality nucleon-nucleon (NN) and three-nucleon (3N) interactions beyond $A=5$ had hindered the *ab initio* approach in the past. Nevertheless progress has been rapid and it has been possible to extend *ab initio* calculations to ${}^7\text{Be}(p, \gamma){}^8\text{B}$, albeit without 3N forces. Navrátil *et al.* (2011) performed such a calculation using the no-core shell model in the continuum with a nucleon-nucleon interaction derived from chiral EFT at fourth

order. The calculation converged with respect to the model space size and reproduces experimental data for the S-factor despite having virtually no free parameters (though the similarity renormalization group evolution parameter was tuned to reproduce the binding energy of ${}^8\text{B}$). The final value of $S_{17}(0) = 19.4(7)$ eV b agrees with the SF II recommendation and, as will be shown later, with our analysis. More recently, a new approach was introduced to combine experimental measurements with *ab initio* predictions, resulting in an “*ab initio*-informed evaluation” Kravvaris *et al.* (2023) that arrived at a value of $S_{17}(0) = 19.8(3)$ eV b.

A significant theoretical advance was the application of the formalism of halo EFT developed by Bertulani *et al.* (2002) to ${}^7\text{Be}(p, \gamma){}^8\text{B}$ calculations by Zhang *et al.* (2015), Zhang *et al.* (2018), and Higa *et al.* (2022). It treats the incoming p and ${}^7\text{Be}$ as point-like particles and the ${}^8\text{B}$ final state as a bound state of the two. The calculation and the associated errors depend upon the accuracy with which one is able to describe the incoming scattering states, the final bound state and the relevant electromagnetic currents. The relative kinetic or CM energy range $E \lesssim 1$ MeV, corresponding to a relative momentum $p \lesssim 40$ MeV/c, is considered within the domain of applicability of halo EFT. The two cluster physical description of ${}^7\text{Be}(p, \gamma){}^8\text{B}$ is expected to hold at ${}^3\text{He}-\alpha$ relative momenta below a physical cutoff $\Lambda \sim 70$ MeV/c set by the threshold (binding momentum) for breaking the ${}^7\text{Be}$ core into these constituents. Further, at energies above the excitation energy of the first excited state, $E_* = 0.4291$ MeV, including the contribution of the excited core ${}^7\text{Be}^*$ is imperative in the EFT.

Zhang *et al.* (2015) and Zhang *et al.* (2018) include such an excited ${}^7\text{Be}$ contribution. Higa *et al.* (2022) provides a calculation without the ${}^7\text{Be}^*$ contribution below E_* and one with it above E_* that also includes an $M1$ contribution to the 1^+ resonance. The two groups’ halo EFT calculations with the ${}^7\text{Be}^*$ contributions differ in several respects that are elaborated in Higa *et al.* (2022). Differences in the final bound state calculation affect the treatment of the short-range interactions and divergences in the EFTs. The interpretation of the asymptotic normalization constants (ANCs) in terms of elastic scattering parameters is affected. However, the momentum dependence of the capture cross section is not impacted by the ANCs, which are fitted to capture data. Thus there is no effect on the cross section. In the incoming channel, Higa *et al.* (2022) use a more general form of the s -wave short-ranged interaction involving the ${}^7\text{Be}^*$ core based on the low-energy symmetry that is included only at NNLO. In contrast, Zhang *et al.* (2015) and Zhang *et al.* (2018) include a short-ranged interaction at LO that could make a difference in the momentum dependence of the cross section. However, as shown in an order-by-order calculation (Higa *et al.*, 2022), the short-ranged interaction is a sub-leading effect so the two halo EFTs should have

similar low-momentum dependence. The third important difference is in the treatment of two-body currents, which affects EFT error estimates. Zhang *et al.* (2015) and Zhang *et al.* (2018) include this at NLO whereas Higa *et al.* (2022) estimate it to be an $N^3\text{LO}$ effect. Bayesian estimates of the two-body current (Zhang *et al.*, 2015, 2018) are not strongly constrained by capture data, consistent with this being a higher-order contribution.

Another widely used approach for describing the cross sections of low-energy nuclear reactions is R-matrix theory. In R-matrix theory the ${}^7\text{Be}(p, \gamma){}^8\text{B}$ cross section can be parametrized in terms of a few parameters representing either “real” or “background” poles. The real poles correspond to the observed states and resonances, while the background terms correspond to the “mean-field” effects that are of the non-resonant type. In the present case for example, a background pole in the 1^- channel is needed so as to reproduce the non-resonant $E1$ part of the capture. The parameters of the real pole corresponding to the bound state are determined by the binding energy and the two ANCs, while the 1^+ resonance parameters are determined by its energy and partial decay widths (both proton and radiative). The R-matrix analysis of this reaction carried out by Barker (1995) arrived at a value of $S_{17}(0) = 17(3)$ eV b; however, it was based on data that were excluded in SF II and here, with the exception of those of Filippone *et al.* (1983).

We fit data using R-matrix theory and halo EFTs, with both frequentist and Bayesian approaches, investigating the impact of different choices for the energy range in which data were fitted, and obtained consistent results. Our recommended $S_{17}(0)$ value and its uncertainty are based on a Bayesian fit of both direct and Coulomb dissociation data taken at energies $E \leq 1250$ keV using the Higa *et al.* (2022) halo EFT $_{*}$. This theory provides a good description of capture data over a wide energy range including the $M1$ contribution from the 1^+ resonance. The theoretical uncertainties are well understood. Further, the contributions of the $S = 1$ and $S = 2$ spin channels to the capture cross section are self-consistently parametrized with the elastic scattering information contained in the known ANCs for this system.

The Bayesian formalism naturally incorporates prior information concerning the ANCs in the fits. The upper limit on the energy in the fit was determined to be high enough to include some data from all the modern experiments, and yet low enough that the contribution of the wide 3^+ resonance at $E \approx 2456$ keV does not exceed 1% of the total cross section at any energy, as estimated by our R-matrix analysis. The fitting procedure is described in the next section, whereas the details, as well as the results of the fits done in different energy ranges and with different theories are reported in the supplemental material.

D. Fitting Procedures

In SF II, it was noted that the different $S_{17}(E)$ data sets had similar energy dependences. The discrepancies among the data sets were primarily due to different absolute normalizations. We address the normalization issue by introducing scaling factors associated with the CMEs. This contrasts with the SF II approach, in which the errors were inflated by the factor $\sqrt{\chi^2/\chi^2(P=0.5)}$. The procedure for handling discrepant data sets adopted here is mathematically rigorous and aligns naturally with the method of estimating theoretical uncertainty discussed below. We prefer this approach to the inflation factor method used in SF II, since the absolute scales of the different data sets can be used as constraints for each other, while there is no information available for a reevaluation of the CMEs. The procedure we used, described below, can be applied to both frequentist and Bayesian analyses.

Although our final recommendation is based on Bayesian analysis alone, both χ^2 minimization (a frequentist approach) and Bayesian posterior probability distribution function evaluation were performed to fit EFTs to the data; in the R-matrix analysis, only the former was used. The likelihood function, defined as the conditional probability $P(D|\boldsymbol{\theta})$ for the data D given theoretical parameters $\boldsymbol{\theta}$, enters both fitting procedures. It can be derived from the χ^2 expression given in D'Agostini (1994)

$$P(D|\boldsymbol{\theta}, \mathbf{s}) = \prod_{\alpha=1}^{N_{\text{set}}} \frac{1}{\sqrt{2\pi\omega_{\alpha}^2}} \exp \left[-\frac{(1-s_{\alpha})^2}{2\omega_{\alpha}^2} \right] \times \prod_{i=1}^{N_{\alpha}} \frac{1}{\sqrt{2\pi\sigma_{\alpha,i}^2}} \exp \left\{ -\frac{[y_{\alpha,i} - \mu_{\alpha,i}(\boldsymbol{\theta})/s_{\alpha}]^2}{2\sigma_{\alpha,i}^2} \right\}, \quad (22)$$

with $\alpha = 1, \dots, N_{\text{set}}$ labeling the different experimental data sets and $i = 1, \dots, N_{\alpha}$ labeling the individual data points of each experiment. In the second term, we divide theoretical predictions $\mu_{\alpha,i}(\boldsymbol{\theta})$ for the data points of the experiment α by a scaling factor s_{α} and consider the combination as our full model for $y_{\alpha,i}$, the i th data point of experiment α with point-to-point uncertainty $\sigma_{\alpha,i}$.

In the first term, we assign s_{α} a Gaussian distribution centered at 1 with a width equal to the CME ω_{α} ; for the GSI, MSU and RIKEN data shown in Table VIII, asymmetric Gaussians are used with ω_{α} depending on the sign of $1 - s_{\alpha}$. In our χ^2 minimization, the χ^2 definition is derived from equating the full exponents in Eq. 22 to $-\chi^2/2$. In the absence of CMEs with $s_{\alpha} = 1$, Eq. 22 leads to the usual uncorrelated least-squares minimization formula.

In the Bayesian analysis, the terms in the first exponential are considered as part of the prior distribution $P(\boldsymbol{\theta}, \mathbf{s})$ that are conditioned on prior knowledge and assumptions such as the CMEs $\boldsymbol{\omega}$ in the scaling factors \mathbf{s} or estimates of theoretical parameters $\boldsymbol{\theta}$ in the EFT. The

posterior distribution for the theoretical parameters is then derived via Bayes's theorem,

$$P(\boldsymbol{\theta}, \mathbf{s}|D) = \frac{P_{\text{Bayes}}(D|\boldsymbol{\theta})P(\boldsymbol{\theta}, \mathbf{s})}{P(D)}, \quad (23)$$

where the likelihood function $P_{\text{Bayes}}(D|\boldsymbol{\theta})$ is defined without the first exponential in Eq. (22). However, since the GSI, MSU and RIKEN data have asymmetric CMEs (see Table VIII), we draw priors from a uniform distribution spanning 50% to 150% of the central values for these that are then used in the asymmetric normalization exponentials in the likelihood function. This was computationally simpler than generating asymmetric Gaussian priors that span the entire $(-\infty, \infty)$ range, which is neither necessary nor physical for a normalization constant.

The constant $P(D)$ in Eq. (23), known as the Bayesian evidence, guarantees the correct normalization of the posterior distribution. It is useful in comparisons of different theories but does not affect the estimation of parameters.

E. Data Analysis and S_{17} Determination

We performed χ^2 fits using several theoretical expressions over multiple energy ranges: $E \leq 475$ keV, $E \leq 1250$ keV, and over a “non-resonant” energy range $E \leq 490$ keV and $805 \text{ keV} \leq E \leq 1250$ keV, in which the $M1$ contribution from the 1^+ resonance can be ignored. Fits with and without Coulomb dissociation data were performed that resulted in overlapping $S_{17}(0)$ determinations within the estimated fitting uncertainties. The different low energy $E \leq 475$ keV fits included one using the halo EFT from Zhang *et al.* (2018) and one using the halo EFT_{*} from Higa *et al.* (2022), both including the excited ${}^7\text{Be}$ contribution. The simpler halo EFT_{gs} expressions, not including an excited component ${}^7\text{Be}^*$, were also fitted in this low energy region with comparable results. For the fits in the other two energy ranges, halo EFT expressions with excited ${}^7\text{Be}^*$ contributions from both Zhang *et al.* (2018) and Higa *et al.* (2022) were used. The Zhang *et al.* (2018) halo EFT expression, which does not include an $M1$ contribution, was supplemented with a Breit-Wigner resonance (Rolfs and Rodney, 1988) with its energy (around the 1^+ resonance energy $E \approx 630$ keV), width, and proton-decay branching ratio determined by the fits. An R-matrix model was fitted to data in the three energy regions with the AZURE2 code (Azuma *et al.*, 2010), again resulting in an $S_{17}(0)$ consistent with the halo EFT values. The results from the various fits are included in the supplemental material.

Bayesian fits were performed with halo EFT expressions from both Zhang *et al.* (2020) and Higa *et al.* (2018) that are compatible with the various χ^2 fits. The Bayesian analysis had advantages over the χ^2 fits in exploring parameter space and in determining realistic uncertainties (King *et al.*, 2019).

The ${}^7\text{Be}(p, \gamma){}^8\text{B}$ capture proceeds through the $S = 2$ and $S = 1$ channels. The reaction is known to be peripheral, resulting in a $S_{17}(0)$ that has a sub-leading dependence on the strong interaction in the initial state (Baye, 2000). This behavior was shown to persist away from the threshold in halo EFT, in which for $E1$ capture the strong interaction in the incoming s -wave channel only contributes at NNLO.

Therefore, to high order, the capture is only sensitive to the sum of the squares of the ANC s in the $S = 2$ and $S = 1$ channels. This results in an arbitrariness in the relative contributions from the two spin channels in the χ^2 fits, without affecting the final S_{17} determination. R-matrix, EFT_{gs}, and EFT_{*} fits confirm this behaviour. On the other hand, the ANCs have been extracted experimentally (Tabacaru *et al.*, 2006; Trache *et al.*, 2003) and also theoretically (Nollett and Wiringa, 2011; Zhang *et al.*, 2018) in *ab initio* calculations. Although the determination of $S_{17}(0)$ is insensitive to the relative contributions of the two spin channels, knowledge of the relative contributions of the ANCs is important in the EFT framework. It establishes the hierarchy of different contributions in the perturbative expansion. This is crucial in developing a self-consistent expansion that is necessary to quantify the theoretical errors at any finite order of the perturbation. Prior knowledge of the ANCs and other parameters can be accounted for naturally in the Bayesian framework in which one specifies the prior probabilities $P(\theta, s)$. Similar constraints can be included in the χ^2 fit by modifying the first exponential in Eq. (22) to center the theoretical parameters around known values.

The Bayesian fits involved six theoretical parameters and 11 normalization constants s_α for the data sets. The priors for the normalization constants were as described earlier. The priors for the theoretical parameters are based on the underlying EFT assumptions (Higa *et al.*, 2022) about the expansion in a small ratio Q/Λ where $Q \sim 40 \text{ MeV}/c$ represents the low momentum of infrared physics and $\Lambda \sim 70 \text{ MeV}/c$ the cutoff of the theory. The 5S_2 and 3S_1 scattering lengths $a_0^{(2)} = -3.18^{+0.55}_{-0.50} \text{ fm}$ and $a_0^{(1)} = 17.34^{+1.11}_{-1.33} \text{ fm}$ (Paneru *et al.*, 2019) are taken as inputs. For $E1$ capture at LO, the only fit parameter, the 5P_2 effective momentum ρ is assumed to scale as Q and accordingly we assume a uniform prior for ρ between $-100 \text{ MeV}/c$ and $1.5 \text{ MeV}/c$, where the upper limit is set by the physical constraint that the ANC squares have to be positive. At NLO, the 3P_2 ANC² $C_{1,1}^2$ is a fit parameter whose prior is taken from a normal distribution determined by experiment. At NNLO, we use the *ab initio* calculation of (Zhang *et al.*, 2018) to draw the ${}^3P_2^*$ ANC² from a normal distribution. A ratio of the s -wave scattering lengths in the coupled-channel 3S_1 - ${}^3S_1^*$ that is assumed to be $\mathcal{O}(1)$ in the EFT is drawn from a normal distribution with a mean of 0 and a standard deviation of 10 to cover a wide range. The $M1$ capture is domi-

nated by the transition ${}^5P_1 \rightarrow {}^5P_2$ and it requires three parameters: a p -wave scattering volume a_1 , an effective momentum r_1 and a two-body current to regulate divergences. Higa *et al.* (2022) determined $a_1 = -108.13 \text{ fm}^3$ and $r_1 = -111.23 \text{ MeV}$ from the narrow 1^+ resonance energy and width. Here, we kept a_1 fixed while drawing r_1 from a uniform distribution between -150 and -50 MeV. The Bayesian fit gave $r_1 \sim -111.23 \text{ MeV}$. The two-body coupling was drawn from a uniform distribution and the fits gave a numerical value consistent with the EFT estimate. We obtain the posterior distribution $P(\theta, s|D)$ using a probabilistic integration method called Nested Sampling (Skilling, 2006) implemented in Python (Feroz *et al.*, 2009) that calculates both the posterior and the evidence.

The experimental data and EFT_{*} fits at LO, NLO, and NNLO are shown in Fig. 8. S_{17} and its derivatives at threshold are given in Table VII. The quality of the fits, the consistency of the numerical values of the fit parameters with EFT assumptions, and the order-by-order improvement in the numerical value of the fitted capture cross section and S-factor give us confidence in the estimated theoretical uncertainty. We find the posterior probability distribution functions to be symmetric and well described by Gaussians, and as such we report the mean and standard deviation in Table VII. Table VIII gives the scaling factors for the experimental data sets determined from the NNLO fit along with their common mode errors.

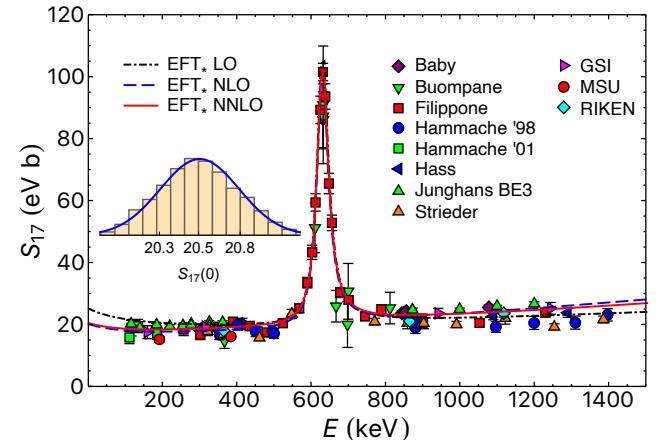


FIG. 8 Astrophysical S-factor for ${}^7\text{Be}(p, \gamma){}^8\text{B}$. Data with $E \leq 1250 \text{ keV}$ were included in the Bayesian fits shown here, which use the halo EFT of Higa *et al.* (2022) that includes an excited ${}^7\text{Be}$ component. The dot-dashed (black) curve is the LO calculation, the dashed (blue) curve is NLO, and the solid (red) curve is NNLO. The inset shows the posterior probability distribution function for $S_{17}(0)$ in eV b. The median as well as the 16th and 84th percentiles of the distribution are specified.

TABLE VII S_{17} and its first two energy derivatives at $E = 50$ eV determined from Bayesian fits of experimental data below 1250 keV using the halo EFT of Higa *et al.* (2022) that includes an excited ${}^7\text{Be}$ component. The first error in each is from the fit. The second error is the estimated LO (30%), NLO (10%), and NNLO (3%) EFT error, respectively, from higher order corrections.

Theory	S_{17} (eV b)	S'_{17}/S_{17} (MeV^{-1})	S''_{17}/S_{17} (MeV^{-2})
LO	25.2(3)(76)	-2.44(4)(73)	35.8(6)(108)
NLO	20.2(2)(20)	-2.02(4)(20)	33.3(6)(33)
NNLO	20.5(3)(6)	-1.86(4)(6)	31.9(6)(10)

TABLE VIII Scaling factors for the data sets at NNLO, the number of data points used in the $E \leq 1250$ keV fits, and the common mode errors.

Data set	Scaling	Data points	CME
Baby	0.97(1)	8	2.22%
Buompane	1.08(3)	9	4.0%
Filippone	1.03(2)	25	11.9%
Hammache '98	1.11(3)	12	5.7%
Hammache '01	1.06(4)	3	11.5%
Hass	1.02(3)	1	3.6%
Junghans BE3	0.93(1)	18	2.67%
Strieder	1.13(2)	9	8.3%
GSI	0.99(3)	4	(-7.7%, +5.6%)
MSU	1.16(4)	2	(-7.1%, +9.6%)
RIKEN	1.05(2)	3	(-12.7%, +8.4%)

F. Theoretical Uncertainty

The theoretical uncertainties in halo EFT are estimated from the perturbative expansion in Q/Λ (Higa *et al.*, 2022). Q is associated with the physics at low momenta of interest: relative momentum $p \lesssim 40$ MeV/ c , binding momentum $\gamma \sim 15$ MeV/ c , inverse Bohr radius $\kappa_C \sim 24$ MeV/ $\hbar c$, excitation momentum $\gamma_\Delta \sim 27$ MeV/ c and resonance momentum $p_R \sim 32$ MeV/ c . The cutoff scale of $\Lambda \sim 70$ MeV is estimated as described above. Except for a narrow momentum range around the 1^+ resonance, capture is dominated by $E1$ transitions from s - and d -waves that preserve the channel spin. Thus the threshold theoretical estimates of $S_{17}(0)$ are based on $E1$ capture. The ANC for the 5P_2 final bound channel is known to be about 5 times larger than the 3P_2 bound channel, so capture in the $S = 2$ channel dominates. The LO contribution in the Q/Λ expansion comes only from s -wave Coulomb interactions in the $S = 2$ channel. The weakly bound ${}^8\text{B}$ state enables peripheral capture without initial state strong interactions.

At NLO, d -wave Coulomb interactions in the $S = 2$ channel and s -wave Coulomb interactions in the $S = 1$ channel are added. Strong interactions in the initial s and d waves of the $S = 1$ channel contribute at NNLO. Perturbatively, the s -wave scattering lengths $a_0^{(1)}$, $a_0^{(2)}$ contribute at this order. Though $a_0^{(1)}$ is much larger than

$a_0^{(2)}$, the $S = 1$ channel is one order higher in the perturbation, making these contributions comparable. We also see mixing in the coupled 3S_1 - ${}^3S_1^*$ channels due to the excited core contribution, which is parameterized by scattering lengths. The mixing requires the ${}^3P_2^*$ ANC in the capture cross section. The theory error is estimated as 30% at LO, 10% at NLO, and 3% at NNLO from an estimated $Q/\Lambda \sim 1/3$. This is validated by the relative sizes of the cross section at different orders of the perturbation, and also by the sizes of the fitted parameters that are consistent with theoretical estimates.

The uncertainty from the higher order contributions is estimated as follows. Two-body currents make a relative contribution $\sim k_0 a_0 L_{E1}$ with photon energy $k_0 = (p^2 + \gamma^2)/(2\mu) \sim Q^3/\Lambda^2$. In the $S = 2$ channel, $a_0^{(2)} \sim 1/\Lambda$ so this is a N^3LO contribution. In the $S = 1$ channel, the larger $a_0^{(1)} \sim 1/Q$ makes this a Q^2/Λ^2 contribution; however, this spin channel is suppressed by one order in the perturbation. Thus two-body currents constitute a N^3LO 3% uncertainty; s -wave effective range corrections would also constitute a N^3LO uncertainty. Initial state d -wave interactions (suppressed by two powers of momentum) in the $S = 2$ channel also constitute a N^3LO uncertainty. The uncertainty in the experimental determination of the s -wave scattering lengths is included in the 3% N^3LO uncertainty.

G. Conclusions and Recommendation for S_{17}

R-matrix and EFT fits of direct and Coulomb dissociation data below 1250 keV, where the influence of resonances other than the 1^+ can be neglected, were carried out. Many different theories and energy ranges were considered and consistent results were obtained. Our recommendation was obtained using the halo EFT of Higa *et al.* (2022) that includes an excited ${}^7\text{Be}$ component. This theory was selected due to the generality and rigor of its treatment of short-range interactions in the initial and final states and of two-body currents, as well as its inclusion of $M1$ transitions and robust uncertainty quantification. The resulting recommendation is $S_{17}(0) = 20.5(3)(6)$ eV b, where the first error is due to fitting and the second due to neglected higher-order contributions; for the purposes of solar modeling, we intend that the two errors be added together in quadrature and recommend $S_{17}(0) = 20.5(7)$ eV b.

The recommended $S_{17}(0)$ has a notably smaller uncertainty than the SF II value for two primary reasons. Significant theoretical developments in halo EFT with a NNLO calculation have reduced the theoretical uncertainty by a factor of two. The use of floating data normalization factors allows discrepant data sets to be brought into agreement for low-energy extrapolation without inflating the measurement uncertainties. In the Bayesian fits, the normalization factors are assigned priors, as is

done for some of the theoretical parameters. Frequentist fits presented in the supplemental material include these in the definition of the χ^2 cost function. Normalization factors different from unity increase the cost, but the overall cost is lowered by bringing precise measurements from different data sets into agreement. This is effective because the number of scale factors is much less than the number of high precision data points. Due to the small uncertainties and large number of data points, the measurement of Junghans *et al.* (2010) exerts a strong influence on the fit. Planned measurements, such as those recently approved at $E = 1$ and 1.2 MeV using a ${}^7\text{Be}$ beam and the DRAGON recoil separator at TRIUMF, must have comparably small statistical and systematic uncertainties in order to compete effectively.

X. The ${}^{14}\text{N}(p, \gamma){}^{15}\text{O}$ reaction ($S_{1\,14}$)

The ${}^{14}\text{N}(p, \gamma){}^{15}\text{O}$ reaction proceeds through a number of bound-state transitions, but the total capture cross section is dominated by the capture to the 6.79 MeV state ($3/2^+$) in ${}^{15}\text{O}$. (In this chapter the reported level energies are taken from the most recent compilation (Ajzenberg-Selove, 1991)). The contribution of this transition to the total reaction cross section at solar energies is 71 – 78% (Adelberger *et al.*, 2011; Azuma *et al.*, 2010). The second strongest transition with a contribution of 16 – 17% is the capture to the ground state of ${}^{15}\text{O}$, while the capture to the 6.18 MeV state ($3/2^-$) is responsible for about 7 – 8% of the total cross section. The remaining 5 – 7% is attributed to the capture to the other bound states (i.e. 5.18 MeV, $1/2^+$; 5.24 MeV, $5/2^+$; 6.86 MeV, $5/2^+$ and 7.28 MeV, $7/2^+$).

In SF II a lower value, with a greatly reduced uncertainty, for $S_{1\,14}(0)$ of 1.66(12) keV b was recommended compared to that of SF I, $3.5^{+0.4}_{-1.6}$ keV b. This was the result of both the new implementation of the more rigorous R-matrix technique (Angulo and Descouvemont, 2001) over the previously used Breit-Wigner analysis, higher precision low-energy capture data sets (Bemmerer *et al.*, 2006; Formicola *et al.*, 2004; Imbriani *et al.*, 2005; Lemut *et al.*, 2006; Marta *et al.*, 2008; Runkle *et al.*, 2005) over the yields presented in Schröder *et al.* (1987), transfer measurements of proton ANC s for the bound states (Bertone *et al.*, 2002; Mukhamedzhanov *et al.*, 2003), and constraints on the lifetime of the 6.79 MeV bound state (Bertone *et al.*, 2001; Schürmann *et al.*, 2008). Since SF II, several new experiments and R-matrix calculations have been completed, but our understanding of the ground state transition still remains incomplete.

A. Current Status and Results

Since SF II, several new measurements have been reported for the capture reaction (Daigle *et al.*, 2016; Frentz *et al.*, 2022; Gyürky *et al.*, 2019b, 2022; Li *et al.*, 2016; Sharma *et al.*, 2020; Wagner *et al.*, 2018), the lifetime of the 6.79 MeV state (Frentz *et al.*, 2021; Galinski *et al.*, 2014; Sharma *et al.*, 2020), the proton bound state ANCs (Artemov *et al.*, 2012), as well as a measurement of the low energy ${}^{14}\text{N}(p, p){}^{14}\text{N}$ scattering cross section (de-Boer *et al.*, 2015). These and past works now include several R-matrix fits (Angulo and Descouvemont, 2001; Formicola *et al.*, 2004; Frentz *et al.*, 2022; Imbriani *et al.*, 2005; Li *et al.*, 2016; Runkle *et al.*, 2005; Wagner *et al.*, 2018), resulting in a range of different approaches to the fitting, which gives insight into the systematic uncertainties of the model.

The recommended partial $S(0)$ are summarised in Table X and discussed in details in sec. X.E. The 6.79 MeV transition, being the strongest, has seen a great deal of experimental attention. Its ease of measurement and relatively well understood reaction contributions result in a small uncertainty in the low energy extrapolation of the S-factor of $\approx 3\%$. The ground state and 6.18 MeV transitions have uncertainties of $\approx 40\%$, owing mainly to the complexity of the underlying reaction contributions and the much smaller cross section.

B. Absolute strength of the 259 keV resonance

The $E = 259$ keV resonance in ${}^{14}\text{N}(p, \gamma){}^{15}\text{O}$ is often used as a reference for normalization in many experiments because it is a narrow resonance, has a precisely known strength, and has γ -ray transitions that span a wide energy range. The absolute strength of the resonance has been determined in three works (Daigle *et al.*, 2016; Gyürky *et al.*, 2019b; Sharma *et al.*, 2020) since SF II. Based largely on the work of Daigle *et al.* (2016), a new resonance strength of $\omega\gamma_{259} = 12.86 \pm 0.45$ meV is recommended, consistent with the previous value but with higher precision thanks to the new data sets. For the averaging we adopted the approach followed by SF II. The common systematic uncertainty of the nitrogen stopping power of 2.9% (Ziegler *et al.*, 2010) was excluded from the weighted mean calculation, and summed in quadrature with the weighted mean uncertainty to obtain the final uncertainty. Additional to this, the stopping power and branching ratio corrected resonance strengths from (Daigle *et al.*, 2016) were used in the calculations instead of the originally quoted values. The resonance strength determination is summarised in Table IX. For the branching ratios, those of Daigle *et al.* (2016) are recommended.

TABLE IX Summary of the $E=259$ keV resonance strength values along with their total uncertainties. The last row gives the recommended value with its total uncertainty.

	$\omega\gamma_{259}$ (meV)
Becker <i>et al.</i> (1982) ^a	13.7 ± 1.0
Runkle <i>et al.</i> (2005) ^a	12.4 ± 0.9
Imbriani <i>et al.</i> (2005)	12.9 ± 0.9
Bemmerer <i>et al.</i> (2006)	12.8 ± 0.6
Daigle <i>et al.</i> (2016)	12.6 ± 0.6
Gyürky <i>et al.</i> (2019b)	13.4 ± 0.8
Sharma <i>et al.</i> (2020)	12.8 ± 0.9
SF III recommended value	12.9 ± 0.5

^a As in (Daigle *et al.*, 2016)

C. Indirect Studies

There are several different indirect techniques that have been used to help constrain the extrapolation of the low energy cross section. A new transfer study of proton ANC s has been reported by Artemov *et al.* (2012). Since the 6.79 MeV level also acts a subthreshold resonance in the ground state transition, its lifetime is also needed. New lifetime measurements have been made by Galinski *et al.* (2014), Sharma *et al.* (2020), and Frentz *et al.* (2021) in addition to the previously available data by (Bertone *et al.*, 2001) and Schürmann *et al.* (2008). Because the lifetime is in a range that is very difficult to access experimentally ($t_{1/2} < 2$ fs) these works still report values either with very large uncertainties or upper limits only. Scattering cross sections can also be very helpful in constraining both bound and unbound level parameters for the R-matrix fit to the capture data. A consistent set of scattering measurements has been made by deBoer *et al.* (2015) superseding the unutilisable data for $^{14}\text{N}(p,p)$ showing a large amount of discrepancy (Gur-bich, 2008). In the new work only yield ratios are reported in order to reduce the uncertainties associated with target degradation. This scattering data give added constraint to the broad, high energy, $3/2^+$ resonance at $E \approx 2.2$ MeV.

D. R-matrix considerations

The phenomenological R-matrix has been the primary model used to fit and extrapolate cross section data, following the analysis of Angulo and Descouvemont (2001). Many of the investigations have used the same multi-channel R-matrix code, AZURE2, which was introduced by Azuma *et al.* (2010). The model has proven to be very popular because it is very flexible in its level structure description, yet at the same time retains enough physical constraints that it can be used for extrapolations into unobserved regions of the cross section. Of course this flexibility also allows for additional complications like over

fitting, mistaking experimental resolution effects for inconsistencies between data sets or the need for additional reaction components (Wiescher *et al.*, 2023), to give just a few common examples.

In the R-matrix framework, the cross section is described by including a number of reaction components, which includes resonances, subthreshold states, and background levels. For radiative capture reactions, an additional component is introduced, external capture (Angulo and Descouvemont, 2001; Barker and Kajino, 1991; Holt *et al.*, 1978), which is used to model the non-resonant capture process. Fig. 9 shows an example multichannel AZURE2 R-matrix fit to scattering data and the three strongest transitions of the $^{14}\text{N}(p,\gamma)^{15}\text{O}$ reaction: 6.79 MeV, ground state, and 6.17 MeV.

Using the alternative R-matrix parameterization of Brune (2002), the resonance components are characterized by observable energies and partial widths, while subthreshold resonances have their strength quantified by the state's asymptotic normalization coefficient (ANC). The strength of the external capture contributions are likewise given by ANCs.

The strongest transition, to the 6.79 MeV state, is dominated by external capture at solar fusion energies. As pointed out in SF II, discrepancies among the data sets over the higher energy range make it unclear whether an additional contribution is needed from a background level. However, as Wagner *et al.* (2018) showed, even when a background level is introduced, it has a very small effect on the low energy extrapolation owing to the dominance of the external capture component.

The next strongest transition is to the ground state, where the R-matrix description of the low energy cross section is rather complicated. It is thought to have significant competing and interfering contributions from the 6.79 MeV subthreshold state, the narrow $E = 259$ keV resonance, external capture, a background level, and the broad, higher energy, resonance at $E = 2.2$ MeV. Despite all of these different contributions, which gives the R-matrix fit a considerable amount of flexibility for this transition, a satisfactory fit to the ground state transition data remains elusive (Frentz *et al.*, 2022; Li *et al.*, 2016; Wagner *et al.*, 2018).

Finally, the 6.18 MeV transition is dominated by the narrow $E = 259$ keV resonance and external capture contributions at low energy. A contribution from the 6.79 MeV subthreshold state may also be significant.

E. Partial S-factors

In this compilation a different approach is adopted to determine the $S(0)$ value for the two strongest transitions than in SF II. Because several are already available in recent publications, no new R-matrix calculation was performed. Instead the reported $S(0)$ values from the

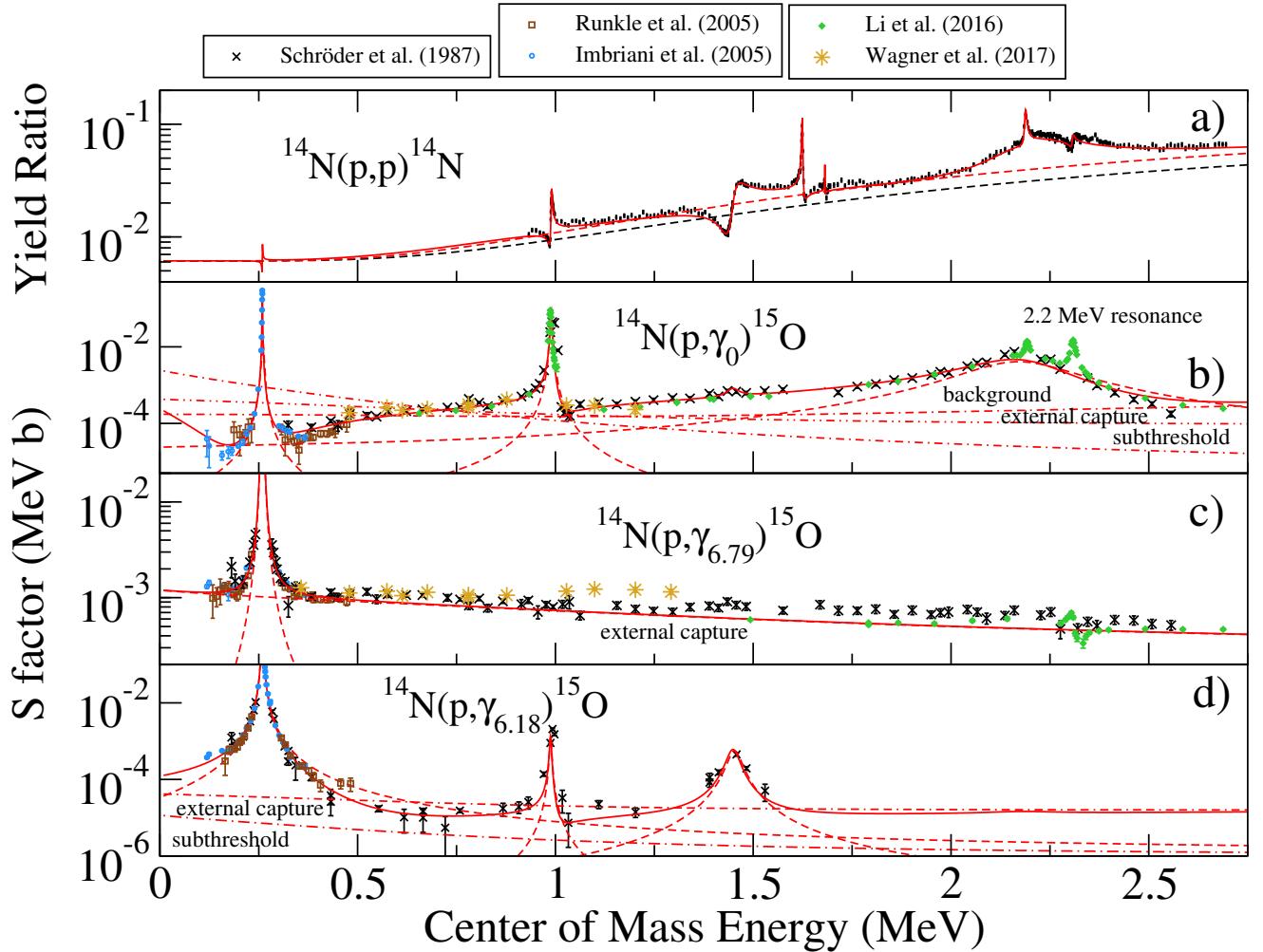


FIG. 9 A representative yield ratio for the $^{14}\text{N}-p$ scattering cross section (deBoer *et al.*, 2015) (a), and low energy S-factors for the three strongest transitions in the $^{14}\text{N}(p,\gamma)^{15}\text{O}$ reaction to the ground (b), 6.79 MeV (c), and 6.18 MeV (d) final states (Imbriani *et al.*, 2005; Li *et al.*, 2016; Runkle *et al.*, 2005; Schröder *et al.*, 1987; Wagner *et al.*, 2018). An example R-matrix fit is shown where individual resonance contributions are indicated by dashed lines, external capture by dotted-dashed lines, and cross sections indicated by solid lines.

original publications (Frentz *et al.*, 2022; Imbriani *et al.*, 2005; Li *et al.*, 2016; Runkle *et al.*, 2005; Wagner *et al.*, 2018) are considered. In the case of Runkle *et al.* (2005), Li *et al.* (2016) and Frentz *et al.* (2022), the data sets are normalized to the strength of the 259 keV resonance, while in Imbriani *et al.* (2005) and Wagner *et al.* (2018) an absolute normalization is established without the resonance strength. For the three former works, the reported $S(0)$ values were re-normalised by the ratio of the resonance strengths used by those works and the present recommended one.

1. Ground State Transition

The ground state transition cross section of the $^{14}\text{N}(p,\gamma)^{15}\text{O}$ reaction is very challenging to measure accurately and to high precision because the cross section

is very low and most experiments perform measurements with high efficiency γ -ray detectors in very close geometry. This increases the yield for these measurements but also greatly increases the effect of γ -ray summing into the ground state transition, with the largest contribution from the 6.79 MeV transition at low energies. In addition, the cross section on the high energy side of the strong, low energy, 259 keV resonance is rather inconsistent between different measurements, likely a result of nitrogen diffusion into the thick target backing material.

In SF II, the R-matrix fit included the data of Schröder *et al.* (1987), Imbriani *et al.* (2005), Runkle *et al.* (2005) and Marta *et al.* (2008), below 2 MeV. A good description of the data was obtained over most of this energy region, with the exception of a consistent over estimation of the experimental data from ≈ 300 to 500 keV. The fit there favored the higher cross sections of Marta *et al.*

(2008), consistently overestimating the data of (Runkle *et al.*, 2005) throughout this energy range. It should also be noted that the data of Schröder *et al.* (1987) were clarified to be yields that were uncorrected for experimental effects. Cross sections were deduced for these data in SF II but were never published.

Since SF II, there have been three new data sets reporting partial cross sections (Frentz *et al.*, 2022; Li *et al.*, 2016; Wagner *et al.*, 2018), which all used the AZURE2 R-matrix code to derive $S(0)$. These values are considered here as an indication of the uncertainty in the low energy extrapolation. Li *et al.* (2016) concentrated on differential cross section measurements between 0.5 and 3.5 MeV, while both Wagner *et al.* (2018) and Frentz *et al.* (2022) focused on bridging the gap between the measurements of Li *et al.* (2016) and the lower energy LUNA (Imbriani *et al.*, 2005) and LENA (Runkle *et al.*, 2005) measurements.

The new data and accompanying R-matrix fits have repeatedly confirmed a discrepancy between the R-matrix fit and the experimental data for the ground state transition, especially when the high energy data of Schröder *et al.* (1987) and Li *et al.* (2016) are included in the fit. On the R-matrix side, this could point to a missing reaction component (i.e. missing resonance or subthreshold state) or an incorrect spin-parity assignment. On the data side, the close geometry setups of most experiments result in large summing corrections. Inaccurate corrections for angular distributions could also play a role. Because of this unresolved inconsistency between data and model, a larger uncertainty has been attributed to this transition compared to SF II (see Table X).

2. 6.79 MeV Transition

In contrast to the ground state, repeated measurements of the 6.79 MeV transition continue to produce consistent results, especially at very low energies. The measurements of Wagner *et al.* (2018) do show a more rapidly increasing cross section at high energies than those of previous works, but are consistent in the low energy region. In addition, the ANC determined from the normalization of the external capture cross section used to model the non-resonant capture in the R-matrix fit yields consistent results with those determined from transfer studies (Bertone *et al.*, 2002; Mukhamedzhanov *et al.*, 2003). Because of this continued consistency across measurements, the value adopted here is the weighted average of the reported values, where the uncertainty has been further reduced from that of SF II.

TABLE X Total $S_{114}(0)$ as the sum of the partial transitions.

Transition	$S_{114}(0)$ (keV b)	$\Delta S_{114}(0)$	Reference
tr \rightarrow 0	0.30 ± 0.11	37%	Present
tr \rightarrow 6.79	1.17 ± 0.03	2.9%	Present
tr \rightarrow 6.17	0.13 ± 0.05	38%	SF II
tr \rightarrow 5.18	0.010 ± 0.003	30%	SF II
tr(5.24) \rightarrow 0	0.068 ± 0.020	30%	SF II
R-matrix sum	$1.68 \pm 0.13^{\text{a}}$	7.6%	
Additional syst. uncert. ^b		3.5%	
Total	1.68 ± 0.14	8.4%	

^a The uncertainty is the sum of the partial transition uncertainties.

^b From the normalisation to the $E = 259$ keV resonance strength.

3. Other Transitions

For the three weaker transitions reported in SF II, there were no new experimental data. Here the values of $S(0)$ have been determined by scaling those from SF II by the ratio of the 259 keV resonance strengths adopted here and from SF II.

F. Total S-factor and conclusions

Figure 9 shows the low energy S-factors and R-matrix fits to the strongest capture transitions and the scattering yields at a representative angle and Table X summarizes the uncertainties estimated for the zero energy S-factor, $S_{114}(0)$ for each of the significant capture transitions. Their sum gives the adopted total $S(0)$ value. The absolute uncertainty of the R-matrix sum is the quadratic sum of the absolute uncertainties of the partial transitions, while the recommended uncertainty conservatively includes an additional 3.5% relative uncertainty from the resonance strength (see Table IX) to which most of the datasets were normalised. As described in Section X.E.1, the uncertainty in the ground state transition has increased to 37% over the 19% estimate of SF II, owing to repeated confirmation of a tension between different experimental data sets and the R-matrix description of the cross section. The uncertainty in the 6.79 MeV transition decreased from 4.0% to 2.9% due to the consistency of the various measurements for this transition.

In addition, the new total capture cross section of Gyürky *et al.* (2022), performed using the activation technique over the energy range from 0.55 to 1.4 MeV, is larger than the sum of the partial R-matrix cross sections by $\approx 25\%$, a significant margin compared to the systematic uncertainties of the different data sets. However, it should be emphasized that for several of the transitions the only data available are those of Schröder *et al.* (1987), which are yields not cross sections and often give only upper limits. For example, the capture to the 7.56 MeV level in this energy range may be significant.

Given the above considerations, a somewhat larger uncertainty of 8.4% has been estimated for $S_{114}(0)$, compared to the 7% uncertainty recommended in SF II. As one of the main sources of uncertainty remains the ground state transition, far geometry angular distribution measurements are recommended extending the differential measurements of Li *et al.* (2016) to lower energies. These measurements will both reduce the effect of summing and provide more information on the underlying reaction mechanisms at these low energies. Further, since the data of Schröder *et al.* (1987) have significant uncorrected experimental effects, new measurements should be made over the wide energy range that they cover to replace them. Such measurements are planned at both the LUNA and JUNA underground facilities. At very low energies, the LUNA and LENA measurements of the ground state transition by Imbriani *et al.* (2005) and Runkle *et al.* (2005), respectively, remain the only measurements. The discrepancy between the two measurements and the tension between the R-matrix fits make new measurements in this range also of high priority. Despite several additional attempts since SF II, the lifetime of the 6.79 MeV transition still remains unmeasured. This remains a high priority, as having a well constrained value will greatly improve our understanding of the low energy behaviour of the ground state transition cross section. Finally, we recommend that the cross section to the 6.18 MeV and 7.56 MeV final states be measured at high energies in order to address the tension between the measured total cross section and the sum of the partial cross sections.

XI. Other CNO, Ne-Na reactions

In the following, we provide information on the proton-induced reactions involving the target nuclei $^{12,13}\text{C}$, ^{15}N , $^{16,17,18}\text{O}$, ^{19}F , $^{20,21,22}\text{Ne}$, and ^{23}Na . Unless otherwise noted, we will focus on the low-energy region, i.e., below 150 keV in the center of mass. The reactions on lighter target nuclei up to and including ^{16}O are of direct relevance for solar neutrinos; the other reactions play a role in hydrogen burning at somewhat higher temperatures, including asymptotic giant branch (AGB) stars.

Table XI lists, where applicable, the S-factor at zero energy, the Gamow peak location, and references. Table XII provides recommended values of the S-factor and corresponding uncertainty on a grid of seven energy values. These results are displayed in Fig. 10, together with their polynomial fits. The fit coefficients are listed in Table I. Table XIII gives values of recommended strengths of narrow resonances. A detailed discussion of each reaction is provided below.

A. Reactions of the CNO-I cycle

1. $^{12}\text{C}(p, \gamma)^{13}\text{N}$

The low-energy reaction cross section is influenced by the tail of a broad resonance and the non-resonant capture to the ground state of ^{13}N . The results of six new direct measurements have been published since SF II (Csedreki *et al.*, 2023; Gyürky *et al.*, 2023; Kettner *et al.*, 2023; Nesaraja *et al.*, 2001; Skowronski *et al.*, 2023a,b).

The first experiment by Nesaraja *et al.* (2001) reports the reaction cross section relative to that of the $^{13}\text{C}(p, \gamma)^{14}\text{N}$ reactions at $E_p^{lab} = 160$ keV to be 0.24 ± 0.03 .

Two of the newest experiments were performed at ATOMKI (Csedreki *et al.*, 2023; Gyürky *et al.*, 2023). The first employed the activation method in a wide laboratory bombarding energy range, $E_p^{lab} = 300 - 1900$ keV, to provide a dataset with an absolute normalization having independent uncertainties, from that of the older data sets obtained using in-beam γ -ray spectroscopy. The main focus of the second experiment was the precise determination of the energies and widths of the two low-energy resonances using in-beam γ -ray spectroscopy with detectors located at several angles.

The fourth experiment, by Skowronski *et al.* (2023b), was performed at the Felsenkeller underground accelerator facility in Dresden, Germany, using in-beam γ -ray spectroscopy. The focus was on determining the parameters of the lowest-lying resonance employing a detector in close geometry with high solid-angle coverage. The data covered the energy range of $E_p^{lab} = 345 - 670$ keV, and allowed for an extrapolation to solar and stellar energies using R-matrix theory. The absolute normalization of this data set is about 25% lower than the previous results. The low-energy slope of the S-factor also changed, resulting in an extrapolated value of $S(25 \text{ keV}) = 1.34 \pm 0.09$ keVb in the solar Gamow peak.

The fifth experiment, by Kettner *et al.* (2023), was performed at Bochum and Notre Dame, and focused on the second resonance in the reaction. Differential cross sections at 0° and 55° , thick target yields, and angular distributions at selected energies were measured in the energy range of $E_p^{lab} = 1000 - 2500$ keV. A comprehensive R-matrix fit of the available data was also performed, resulting in $S(25 \text{ keV}) = 1.48 \pm 0.09$ keVb.

The sixth experiment, by the LUNA collaboration (Skowronski *et al.*, 2023a), obtained data in the range of $E_p = 80 - 400$ keV, reaching the lowest bombarding energies to date. Both the activation technique and in-beam γ -ray spectroscopy were applied, with a consistent absolute normalization of the data, despite the different systematic uncertainties introduced by the two methods. Their S-factor in the solar Gamow peak, $S(25 \text{ keV}) = 1.53 \pm 0.06$ keVb, is in good agreement with Kettner *et al.* (2023), and is compatible within 2σ with

Skowronski *et al.* (2023b).

Apart from the direct measurements discussed above, the ground state ANC was determined in a recent particle transfer experiment (Artemov *et al.*, 2022). That work also reported a fit of the radiative capture data available at the time. The R-matrix fits to the data of Kettner *et al.* (2023); Skowronski *et al.* (2023a,b) give ANC values consistent with the result of the transfer work.

Thanks to the high consistency of all recent datasets, the extrapolation of the cross section to solar energies has now improved uncertainty. Owing to the self-consistency between the used methods and the precise data covering the lowest energies to date we adopt the S-factor recommended by Skowronski *et al.* (2023a) (see Tables XI and XII).

2. $^{13}\text{C}(p, \gamma)^{14}\text{N}$

This reaction proceeds to the ground state and five excited states in ^{14}N . The low-energy cross section is influenced by the tail of a broad resonance at laboratory bombarding energy of $E_r^{\text{lab}} \approx 558$ keV. Previously, only a single work reported cross sections for all transitions (King *et al.*, 1994), which was the basis of the S-factor recommendations in both SF I and SF II.

A new relative measurements was carried out by Nesaraja *et al.* (2001) as mentioned in the previous section.

Later a new comprehensive experiment by the LUNA collaboration (Skowronski *et al.*, 2023a) provided data in the energy range of $E_p^{\text{lab}} = 80 - 400$ keV. One of their data sets was obtained with HPGe detectors to study partial cross sections. Their second data set was obtained with a γ -ray calorimeter to measure total reaction cross sections. The results from both data sets agree in the absolute cross section scale, despite major differences in the systematic uncertainties of the two methods. The new results are about 30% lower than previous results of King *et al.* (1994). Despite this scale difference the calculated S-factor ratio for the $^{12}\text{C}(p, \gamma)^{13}\text{N}$ and $^{13}\text{C}(p, \gamma)^{14}\text{N}$ reactions at $E_p^{\text{lab}} = 160$ keV is 0.26 ± 0.02 , in perfect agreement with the experimental reported one of 0.24 ± 0.03 by Nesaraja *et al.* (2001).

The adopted S-factor given in Tables XI and XII is based on the new result of Skowronski *et al.* (2023a), owing to its self-consistency between the different methods for measuring the cross sections, and the higher precision of the dataset compared to King *et al.* (1994).

3. $^{15}\text{N}(p, \alpha)^{12}\text{C}$

Following SF II, the experimental and phenomenological work on the $^{15}\text{N}(p, \alpha)^{12}\text{C}$ reaction can be divided into three groups addressing different aspects: (i) new (p, α_0) data (deBoer *et al.*, 2012); (ii) new $(p, \alpha\gamma)$ data (de-

Boer *et al.*, 2021a; Gyürky *et al.*, 2019a; Imbriani *et al.*, 2012a,b; Reinhardt *et al.*, 2016); and (iii) R-matrix analysis of existing data sets (Brune and deBoer, 2020; deBoer *et al.*, 2013).

Other studies have been conducted for laboratory bombarding energies above 400 keV. In Reinhardt *et al.* (2016), the strength of a resonance at $E_r^{\text{c.m.}} \approx 403$ keV (2^- , $E_x = 12530$ keV) is tabulated, and the value is in agreement with previous determinations. In Gyürky *et al.* (2019a), the thick-target yield for the $E_r^{\text{c.m.}} = 841$ keV resonance is shown, but no further details are reported. New cross section data for the $(p, \alpha_1\gamma)$ reaction down to about 900 keV center-of-mass energy are displayed in deBoer *et al.* (2021a). They used the data for a comprehensive R-matrix fit, and reported parameters of resonances above about 1500 keV center-of-mass energy.

Imbriani *et al.* (2012a,b) measured the $^{15}\text{N}(p, \alpha\gamma)^{12}\text{C}$ reaction down to about 140 keV laboratory bombarding energy, and showed that the S-factor of the $(p, \alpha_1\gamma)$ channel is orders of magnitude lower compared to the (p, α_0) channel.

Finally, deBoer *et al.* (2013) contains a detailed analysis of all available channels populating the ^{16}O compound nucleus, expanding the results presented in deBoer *et al.* (2012). In this paper, a wide range of experimental data were considered leading to a single consistent R-matrix fit, accurately describing the broad level structure of ^{16}O below $E_x = 13500$ keV. The resulting fit was used to extract an improved determination of the low-energy S-factor for the $^{15}\text{N}(p, \alpha)^{12}\text{C}$ reaction. In addition, Brune and deBoer (2020) have developed the partial-wave formalism for calculating the angular distributions of secondary γ -rays following particle emission, providing the framework needed for future studies that can directly utilize differential measurements of the $^{15}\text{N}(p, \alpha_1\gamma)^{12}\text{C}$ reaction.

The available evidence confirms that the (p, α_0) channel dominates the total S-factor. The global R-matrix analysis in deBoer *et al.* (2013) provides results that slightly exceed the experimental data of Redder *et al.* (1982) and also estimates a slightly larger value, $S(0) = (95 \pm 6) \times 10^3$ keVb, than previous works (see Table VII of deBoer *et al.* (2013)), calling for additional low-energy measurements. The extrapolated S-factor in deBoer *et al.* (2013) agrees with the findings in La Cognata *et al.* (2009), which is the most recent indirect study, determining the S-factor down to 20 keV center-of-mass energy. The global R-matrix analysis suggests a renormalization of the La Cognata *et al.* (2009) S-factor by 6%, which is within the quoted uncertainty. Based on the information provided above, we adopt the S-factor of La Cognata *et al.* (2009), $S(0) = (73 \pm 5) \times 10^3$ keVb, which was also adopted in SF II.

B. Reactions of the NO, or CNO-II, cycle

1. $^{15}\text{N}(p, \gamma)^{16}\text{O}$

The low-energy $^{15}\text{N}(p, \gamma)^{16}\text{O}$ cross section is dominated by its ground state transition (Imbriani *et al.*, 2012b; Leblanc *et al.*, 2010). This transition's cross section is enhanced by the constructively interfering tails of two broad resonances at laboratory proton energies of $E^{lab} = 338$ and 1028 keV, which correspond to 1^- levels in the ^{16}O compound system. In addition to ground-state proton and γ -ray de-excitation, these levels also decay strongly through the α_0 and α_1 channels. Currently, the favored framework for the analysis and extrapolation of the available data is the phenomenological R-matrix. Here, the main components that are used to model the low-energy cross section of the $^{15}\text{N}(p, \gamma)^{16}\text{O}$ reaction are the two broad 1^- resonances, external capture, and a 1^- background level. The strength of the external capture can be characterized by the ground state proton ANC of ^{16}O , which has been experimentally determined via a proton transfer reaction (Mukhamedzhanov *et al.*, 2008). Gamma-ray de-excitation to the ground state has also been observed through a 2^+ level in this energy region in the $^{12}\text{C}+\alpha$ reactions, but has been shown to be negligibly weak for $^{15}\text{N}(p, \gamma)^{16}\text{O}$ (deBoer *et al.*, 2013).

The cross section measurements used in SF II were those of Hebbard (1960), Rolfs and Rodney (1974), and Bemmerer *et al.* (2009). While the then newly measured data of Bemmerer *et al.* (2009) represented a significant improvement in accuracy over Rolfs and Rodney (1974) and in precision over Hebbard (1960), they only cover a limited energy range. For SF II, the two-level R-matrix fit of Mukhamedzhanov *et al.* (2008) was adopted, which simultaneously fit $^{15}\text{N}(p, \alpha)^{12}\text{C}$ data, giving $S(0) = 36 \pm 6$ keVb for the $^{15}\text{N}(p, \gamma)^{16}\text{O}$ reaction (see also Barker (2008)). Importantly, it showed that the extrapolated $S(0)$ value of Rolfs and Rodney (1974), 64 ± 6 keVb, was an overestimate.

Following SF II, a comprehensive measurement campaign was carried out at LUNA and the University of Notre Dame (Leblanc *et al.*, 2010), covering a laboratory energy range from $E^{lab} = 130$ to 1800 keV. In that work, a three-level R-matrix fit was used, but the ground state proton ANC of Mukhamedzhanov *et al.* (2008) was not considered, resulting in a somewhat larger extrapolated value of $S(0) = 39.6(26)$ keVb than that adopted in SF II. Mukhamedzhanov *et al.* (2011) subsequently showed that a three-level R-matrix fit of equal quality could be obtained with the ANC, and reported an extrapolated value of $S(0)$ that ranged from 33 to 40 keVb.

Caciolli *et al.* (2011) expanded the energy range of Bemmerer *et al.* (2009) by measuring the cross section up to $E_{c.m.} = 370$ keV. Their experiment scanned over the top of the lowest-energy broad resonance at $E_{c.m.} = 312$ keV. The data were found to be consistent with the ground

state data of Leblanc *et al.* (2010).

deBoer *et al.* (2013) performed a comprehensive R-matrix fit that included not only the $^{15}\text{N}+p$ data, but also a wide range of $^{12}\text{C}+\alpha$ data over a similar excitation energy range (see also Section XI.A.3). The value from this analysis, $S(0) = 40 \pm 3$ keVb, was slightly above that of Leblanc *et al.* (2010) and agreed with the upper limit of Mukhamedzhanov *et al.* (2011). However, it is lower than the recent potential model extrapolation of $S(0) = 45^{+9}_{-7}$ keVb found in the NACRE II compilation (Xu *et al.*, 2013). The difference between the two fits is likely due to the fact that NACRE II disregarded the higher-energy part of the Leblanc *et al.* (2010) data, which maps the interference pattern between the two 1^- resonances more accurately than older data sets. In order to reproduce experimental data in that region, a 1^- background level was added in the R-matrix fits of Leblanc *et al.* (2010) and later works.

Here we adopt the value of $S(0) = 40 \pm 3$ keVb from deBoer *et al.* (2013) as it considers the most comprehensive set of available data.

While low-energy future measurements are recommended, experiments at high energy are highly desirable to better understand the reaction mechanisms responsible for the interference pattern that is observed between the two well known 1^- resonances, which can not be completely reproduced unless some background contribution is also included.

2. $^{16}\text{O}(p, \gamma)^{17}\text{F}$

Below a laboratory energy of about 2700 keV, i.e., the energy of the lowest-lying resonance (Tilley *et al.*, 1993), the $^{16}\text{O}(p, \gamma)^{17}\text{F}$ reaction is a prime example of the non-resonant direct radiative capture process. This reaction has been measured many times using a variety of techniques, including activation, prompt γ -ray detection, and inverse-kinematics experiments. The $E1$ capture can proceed to the ground or first-excited state in ^{17}F , with the latter being the dominant transition. In SF II, a value of $S(0) = 10.6 \pm 0.8$ keVb with a 7.5% uncertainty was adopted. This result was obtained by normalizing the microscopic-model calculations of Baye *et al.* (1998) to the data of Rolfs (1973) and Morlock *et al.* (1997).

A recent statistically rigorous evaluation of the S-factor data of Becker *et al.* (1982); Chow *et al.* (1975); Hester *et al.* (1958); Morlock *et al.* (1997) is presented in Iliadis *et al.* (2022). Their analysis, including a discussion of experimental uncertainties, used a combined fit of the transitions to the ground and first-excited states in ^{17}F , and their sum, and was performed using a Bayesian model. The physical model was a single-particle model employing a Woods-Saxon potential for generating the radial bound-state wave function. The fit had three adjustable parameters: the radius parameter and diffuse-

ness of the Woods-Saxon potential, and the ANCs for scaling the theoretical non-resonant capture S-factors. It was also found that a poor fit was obtained when using spectroscopic factors instead as scaling parameters. As the $^{16}\text{O}(p, \gamma)^{17}\text{F}$ reaction at low bombarding energies is peripheral, the analysis of the S-factor in terms of ANCs greatly reduced the sensitivity of the fit to the single-particle potential parameters. For the ANCs of the ground and first-excited state transitions, they found values of $C_{\text{gs}}^2 = 1.115 \text{ fm}^{-1} (\pm 4.0\%)$ and $C_{fes}^2 = 7063 \text{ fm}^{-1} (\pm 4.0\%)$, respectively. These results agree with those measured in the $^{16}\text{O}(^{3}\text{He}, \text{d})^{17}\text{F}$ reaction by Gagliardi *et al.* (1999). It is not surprising that the uncertainties ($\approx 10\%$) in the latter reference are significantly larger than for the values quoted above, because the distorted wave Born approximation (DWBA) analysis of Gagliardi *et al.* (1999) was subject to ambiguities in the choice of optical model potentials for the incoming and outgoing channels.

The total S-factor at zero energy reported by Iliadis *et al.* (2022) was $S_{\text{tot}}(0) = 10.92 \text{ keVb} (\pm 4.0\%)$. The uncertainty is significantly smaller compared to the results published by Iliadis *et al.* (2008), which were derived from an analysis that was not statistically rigorous. We adopt the results of Iliadis *et al.* (2022) in this work. Their S-factor is presented in Table XII.

We endorse the two recommendations made by Iliadis *et al.* (2022). First, they advocated for a new low-energy measurement of the $^{16}\text{O}(p, \gamma)^{17}\text{F}$ reaction, specifically at center-of-mass energies below 200 keV, which have so far only been reached in the experiment of Hester *et al.* (1958), albeit with a relatively large systematic uncertainty of 14%. Second, it would be interesting to compare their results to those from a future R-matrix analysis of the same data sets.

3. $^{17}\text{O}(p, \gamma)^{18}\text{F}$

In SF II, only the non-resonant component of the $^{17}\text{O}(p, \gamma)^{18}\text{F}$ reaction cross section was discussed, which dominates the total cross section at solar temperatures. At the time, the results of Chafa *et al.* (2007) were recommended. Since the publication of SF II, several new S-factor measurements have been reported, both at low and high energies (Buckner *et al.*, 2015; Di Leva *et al.*, 2014; Gyürky *et al.*, 2017; Hager *et al.*, 2012; Kontos *et al.*, 2012; Newton *et al.*, 2010; Scott *et al.*, 2012). Despite differences in the techniques exploited and energy ranges covered, the derived $S(0)$ values agree across different experiments, both for the non-resonant capture and the total S-factor, where the latter quantity includes the low-energy tails of two broad resonances ($E_r^{\text{c.m.}} = 557$ and 677 keV). Here we recommend the total S-factor from the most recent analysis of (Gyürky *et al.*, 2017). The $^{17}\text{O}(p, \gamma)^{18}\text{F}$ reaction cross section was measured in

the laboratory bombarding energy range of 500–1800 keV with a typical total uncertainty of 10% using the activation method. A comparison with other independent results, when possible, at $E_p^{\text{lab}} = 500 \text{ keV}$ resulted in good agreement with Hager *et al.* (2012) and Kontos *et al.* (2012), while a 20% difference is observed with respect to Newton *et al.* (2010). An R-matrix analysis was performed considering all the low energy data, resulting in a value of $S(0)_{\text{tot}} = (4.7 \pm 1.0) \text{ keVb}$ ⁸.

At temperatures near 60 MK, typical for hydrogen-shell burning in AGB stars, the $^{17}\text{O}(p, \gamma)^{18}\text{F}$ reaction rate is dominated by the $E_r^{\text{c.m.}} = 65 \text{ keV}$ resonance, which was not considered in SF II. No direct measurements are reported for the resonance strength, $\omega\gamma$, which is presently estimated from experimental values of the partial and total widths. Partial widths of $\Gamma_\alpha = (130 \pm 5) \text{ eV}$ and $\Gamma_\gamma = (0.44 \pm 0.02) \text{ eV}$ were measured using the $^{14}\text{N} + \alpha$ reaction (Berka *et al.*, 1977; Mak *et al.*, 1980). The most uncertain quantity is the proton partial width, Γ_p , which is estimated from the measured strength of the corresponding resonance in the (p, α) channel. Our best estimate for the (p, α) strength is $\omega\gamma_{p\alpha} = 4.6_{-1.2}^{+5.4} \times 10^{-9} \text{ eV}$ (see Table XIII and the following section for details), which leads to $\Gamma_p = 18.5_{-4.7}^{+21.7} \times 10^{-9} \text{ eV}$. Hence, our recommended value is $\omega\gamma_{p\gamma} = 15.6_{-4.0}^{+18.3} \times 10^{-12} \text{ eV}$ (see Table XIII). This value is consistent with the strength recommended by Palmerini *et al.* (2013), though we have inflated the error to account for the uncertain status of the proton width. A direct measurement of the (p, γ) strength of the $E_r^{\text{c.m.}} = 65 \text{ keV}$ resonance is ongoing at LUNA (Ciani *et al.*, 2022). Moreover, the 65 keV (1^-) resonance is expected to interfere with the –2.0 keV (1^-) subthreshold resonance. We recommend the treatment of the interference contribution described in Buckner *et al.* (2015). At typical temperatures of classical novae, the main contributor to the $^{17}\text{O}(p, \gamma)^{18}\text{F}$ reaction rate is the $E_r^{\text{c.m.}} = 183.9 \text{ keV}$ resonance, which was first observed by Fox *et al.* (2004). It was subsequently remeasured by Buckner *et al.* (2015); Chafa *et al.* (2007); Di Leva *et al.* (2014); Scott *et al.* (2012). Here we recommend the literature-data weighted mean adopted in Buckner *et al.* (2015), $\omega\gamma = (1.77 \pm 0.09) \times 10^{-6} \text{ eV}$.⁹

4. $^{17}\text{O}(p, \alpha)^{14}\text{N}$

The $^{17}\text{O}(p, \alpha)^{14}\text{N}$ reaction rate is determined by the contributions of narrow resonances. At solar energies,

⁸ In the publication of Gyürky *et al.* (2017) the uncertainty is referred to as statistical only. However, it represents, in fact, the total uncertainty (Gy. Gyürky, priv. comm.)

⁹ The article by Gesuè *et al.* (2024) reports an $\omega\gamma$ bare value of $(30 \pm 6(\text{stat}) \pm 2(\text{syst})) \times 10^{-12} \text{ eV}$ for the 64.5 keV resonance strength.

the dominant contributions originate from resonances at $E_r^{c.m.} = -2$ and 65 keV. At typical temperatures of AGB stars and classical novae, the main contributor to the rate is a resonance at $E_r^{c.m.} = 183.9$ keV. Concerning the subthreshold resonance, no new results have been published since SF II.

The $E_r^{c.m.} = 65$ -keV resonance has been studied recently both in a direct experiment ($\omega\gamma = (10.0 \pm 1.4_{stat} \pm 0.7_{syst}) \times 10^{-9}$ eV; Bruno et al. 2016) and with the Trojan Horse indirect Method ($\omega\gamma = (3.42 \pm 0.60) \times 10^{-9}$ eV; Sergi et al. 2015). The resonance strength reported by Bruno et al. (2016) deviates by a factor of 2.5 from that of Sergi et al. (2015) and a factor of 2 from the pre-SF II value of Fox et al. (2005), which includes a revised analysis of the data from Blackmon et al. (1995). Since the reason behind the discrepancies between literature strengths remains unclear, an independent investigation of this resonance would be desirable. Here we recommend a weighted average of the available literature strengths, $\omega\gamma = 4.6_{-1.2}^{+5.4} \times 10^{-9}$ eV. The uncertainty was enlarged to include all literature values.

The resonance at $E_r^{c.m.} = 183.9$ keV, on the other hand, is well known. The most recent determination of its strength is reported in Bruno et al. (2015), i.e., $\omega\gamma = (1.68 \pm 0.03_{stat} \pm 0.12_{syst}) \times 10^{-3}$ eV, which is in excellent agreement with previous results (Chafa et al., 2007; Moazen et al., 2007; Newton et al., 2007).

C. Reactions of the CNO-III cycle

1. $^{18}\text{O}(p, \gamma)^{19}\text{F}$

Important components for this reaction channel are a $E_r^{c.m.} = 20$ keV resonance, for which only indirect data exist (Champagne and Pitt, 1986; La Cognata et al., 2008; Wiescher et al., 1980), the resonances at $E_r^{c.m.} = 90$ keV and $E_r^{c.m.} = 143$ keV and the non-resonant capture process. Reported strengths of the $E_r^{c.m.} = 20$ keV resonance vary by one order of magnitude between 1.4×10^{-22} (Lorenz-Wirzba et al., 1979) and 1.9×10^{-21} eV (Wiescher et al., 1980), depending on the assumed proton widths. We recommend a value of $\omega\gamma = 5.7 \times 10^{-22}$ eV, based on the results of Champagne and Pitt (1986). A comprehensive measurement was done by Wiescher et al. (1980), covering the 80–2200 keV range at multiple angles and measuring branching ratios, the direct component (extrapolated from high-energy data), and strengths down to the $E_r^{c.m.} = 143$ keV resonance. The strength of the latter resonance, $\omega\gamma = (0.98 \pm 0.03) \times 10^{-3}$ eV, has been remeasured multiple times with consistent results (Becker et al., 1982; Best et al., 2019; Dermigny et al., 2016; Pantaleo et al., 2021; Vogelaar et al., 1990; Wiescher et al., 1980). The strength of the $E_r^{c.m.} = 90$ keV resonance was recently controversial (Buckner et al., 2012; Fortune, 2013), but a measure-

ment by Best et al. (2019) found it to be of insignificant strength ($\omega\gamma = (0.5 \pm 0.1) \times 10^{-9}$ eV). The non-resonant component was determined in Wiescher et al. at an energy of $E^{c.m.} = 1752$ keV and extrapolated to a value of $S(0) = 15.7$ keVb. Buckner et al. normalized a non-resonant capture model calculation to the experimental value at $E^{c.m.} = 1752$ keV, resulting in $S(0) = 7.1$ keVb. No uncertainties were given for these two S-factors. Best et al. (2019) extracted the non-resonant component from a cross section fit at center-of-mass energies between 85 and 150 keV, resulting in $S(0) = (23 \pm 3.8)$ keVb. Here we adopt this value, but further investigation of the low-energy cross section is necessary to resolve these discrepancies.

2. $^{18}\text{O}(p, \alpha)^{15}\text{N}$

The low-energy trend of the $^{18}\text{O}(p, \alpha)^{15}\text{N}$ cross section is determined by three resonances at 20, 90 and 143 keV (see Table XIII). The most recent determination of the strength of the 20 keV resonance comes from a THM experiment (La Cognata et al., 2008). Their result was already recommended in SF II and we maintain the same recommendation. The resonance at 90 keV was investigated in the past with direct (Lorenz-Wirzba et al., 1979) and indirect (La Cognata et al., 2008) techniques, finding consistent results. Nevertheless, a new direct measurement reported a resonance strength one order of magnitude higher than the previously adopted value (Bruno et al., 2019). Since no clear explanation can be found for such a discrepancy, here we adopt a weighted average of the three resonance strengths, but we inflate the uncertainty so that it embraces all three values. New experiments to solve this discrepancy are advised. The resonance at 143 keV is well known, and different experiments provided consistent results (Becker et al., 1995; Bruno et al., 2015; Lorenz-Wirzba et al., 1979). Here we adopt a weighted average of Becker et al. (1995) and Bruno et al. (2015), which have independent systematic uncertainties (while no information on individual contributions to the total uncertainty is given in Lorenz-Wirzba et al. (1979)). The non-resonant differential cross section has recently been measured down to $E_{c.m.} = 55$ keV by Bruno et al. (2019). At overlapping energies, the S-factor is consistent with the previous measurement by Lorenz-Wirzba et al. (1979). Bruno et al. (2019) found a broad structure at $E \approx 110$ keV that they could only explain assuming the existence of a so-far unobserved resonance at $E_r^{c.m.} = 106$ keV. Further investigations are needed to shed light on this supposed resonance.

3. $^{19}\text{F}(p, \gamma)^{20}\text{Ne}$

The $^{19}\text{F}(p, \gamma)^{20}\text{Ne}$ reaction has been challenging to access experimentally because it has a much smaller cross section than the competing $^{19}\text{F}(p, \alpha\gamma)^{16}\text{O}$ reaction. The cross section of both reactions is dominated by a mixture of narrow and broad resonances corresponding to mostly unnatural parity states in ^{20}Ne (mainly $J^\pi = 1^+$ and 2^-). At very low energies, the cross section may be enhanced by a near threshold state at $E_{\text{c.m.}} \approx 11$ keV and subthreshold and non-resonant capture contributions may also contribute significantly, but have not yet been well characterized (deBoer *et al.*, 2021b), as few particle transfer studies have been published (Betts *et al.*, 1975; Kious, 1990).

The first low-energy cross section measurement was that of Couture *et al.* (2008), who measured only the $^{19}\text{F}(p, \gamma_1)^{20}\text{Ne}$ transition. While this was likely a strong transition, there can certainly be other transitions with sizable contributions. Couture *et al.* (2008) covered the center-of-mass energy range from 200 to 760 keV, where three strong resonances were observed at $E_r^{\text{c.m.}} = 323.9$, 564 and 634 keV. Only an upper limit was reported for an expected resonance at $E_r^{\text{c.m.}} = 213$ keV. Also, no evidence of either a near-threshold state, subthreshold state, or non-resonant capture has been observed.

An investigation of the γ -ray branching ratios of the $E_r^{\text{c.m.}} = 323.9$ keV resonance was undertaken by Williams *et al.* (2021), using the DRAGON recoil separator at TRIUMF. Recoil detection provided an effective method of discriminating the large background from the $^{19}\text{F}(p, \alpha\gamma)^{16}\text{O}$ reaction. It was found that the primary transitions to both the ground and the 4967 keV state dominate over the primary transition to the 1633 keV first-excited-state.

Finally, a new direct measurement of the low-energy cross section has been reported by Zhang *et al.* (2022), which overlaps with the previous low-energy data of Couture *et al.* (2008). A new resonance was discovered at $E_r^{\text{c.m.}} = 225.2$ keV in the $^{19}\text{F}(p, \gamma_1)^{20}\text{Ne}$ cross section, attributable to a 3^- state that corresponds to a weak resonance observed in the $^{19}\text{F}(p, \alpha\gamma)^{16}\text{O}$ data of Spyrou *et al.* (2000) at the same energy. This low-energy resonance increases the $^{19}\text{F}(p, \gamma)^{20}\text{Ne}$ reaction rate by about an order of magnitude at $T \approx 0.1$ GK. Like Williams *et al.* (2021), Zhang *et al.* (2022) also observed a sizeable ground state transition for the $E_r^{\text{c.m.}} = 323.9$ keV resonance. As in the $^{19}\text{F}(p, \alpha\gamma)^{16}\text{O}$ reaction (see Section XI.C.4), an upper limit was also estimated for the proposed $E_r^{\text{c.m.}} = 11$ keV resonance.

We have therefore adopted the resonance energies and strengths of Zhang *et al.* (2022), which are in good agreement with the previous measurements of Couture *et al.* (2008) and Williams *et al.* (2021) in the overlapping energy region (see Tables XI and XIII).

4. $^{19}\text{F}(p, \alpha)^{16}\text{O}$

The $^{19}\text{F}(p, \alpha)^{16}\text{O}$ reaction was not included in SF II. NACRE (Angulo *et al.*, 1999) pointed out that the total astrophysical S-factor was dominated by the non-resonant (p, α_0) channel, though underscoring the large uncertainties ($\approx 50\%$) affecting the reaction rate at low temperatures. After NACRE, the situation greatly changed following the publication of a large number of papers addressing the measurement of the α_0 channel (Indelicato *et al.*, 2017; La Cognata *et al.*, 2011, 2015a; Lombardo *et al.*, 2013, 2015), of the $\alpha\gamma$ channel (Ding *et al.*, 2002; Fessler *et al.*, 2000; Pham *et al.*, 2021; Spyrou *et al.*, 2000; Zhang *et al.*, 2019, 2021b). In addition, several authors attempted to reach a consensus astrophysical factor carrying out refined analyses using, for instance, the phenomenological R-matrix (deBoer *et al.*, 2021b; He *et al.*, 2018; Lombardo, 2019; Lombardo *et al.*, 2019; Zhang *et al.*, 2021a), and carried out new theoretical calculations of the astrophysical S-factor (Sahoo and Basu, 2021).

Regarding the measurement of the astrophysical factor for the α_0 channel, populating the ^{16}O ground state, indirect (using the THM (Tribble *et al.*, 2014)) and direct methods were used, leading to consistent results. While the direct measurement (Lombardo *et al.*, 2015) could reach about 170 keV in the center-of-mass, the THM measurements (Indelicato *et al.*, 2017; La Cognata *et al.*, 2011, 2015a) could reach 5 keV, providing an extrapolated $S(0) = (17.4 \pm 2.7) \times 10^3$ keVb. The most striking feature of both direct and indirect measurements is the evidence of resonances populating the $E_{\text{c.m.}} \lesssim 600$ keV energy region, at odds with the NACRE extrapolation.

Until Zhang *et al.* (2021b), direct measurements of the $\alpha\gamma$ channel stopped at energies well above the Gamow energy. Zhang *et al.* (2021b) could reach energies as low as ≈ 72 keV in the center of mass, still above the Gamow window for solar fusion, thanks to the use of the Jinping Underground Nuclear Astrophysics experimental facility (JUNA). Though the lowest energy point is affected by a large uncertainty, 87%, and no correction for the electron screening effect is reported (for instance, Assenbaum *et al.* (1987) assume a 20% enhancement at 70 keV center-of-mass energy for p -induced reactions on oxygen), the R-matrix analysis the authors carried out suggests the occurrence of a $E_r^{\text{c.m.}} = 11$ keV 1^+ level at 12855 keV ^{20}Ne excitation energy, which might significantly enhance the astrophysical factor, leading to a $S(0) \approx 200$ MeVb. If this result is confirmed, the $\alpha\gamma$ channel will dominate the low-energy total astrophysical S-factor (below about 50 keV in the center of mass) of the $^{19}\text{F}(p, \alpha)^{16}\text{O}$ reaction. If the $E_r^{\text{c.m.}} = 11$ keV resonance is excluded from the fit, $S(0) \approx 1$ MeVb would be deduced.

The presence of new data sets triggered the publication of review papers aiming to provide the best fit of the

available data. In particular, Lombardo *et al.* (2019) carried out a comprehensive R-matrix analysis spanning an energy range from about $E^{c.m.} = 200$ keV to 12000 keV. The R-matrix fit predicts a $S(0) \approx 10$ MeVb for the α_0 channel, with a conservative uncertainty of about 20%, lower than the value from Indelicato *et al.* (2017). An additional interesting result from this work is that the α_1 channel, corresponding to the population of the 0^+ ^{16}O first excited state, is negligible with respect to the α_0 one below about $E^{c.m.} = 50$ keV, yet if the 2^+ state at 13095 keV is assumed to contribute, as supported by the $^{16}\text{O}(\alpha, \alpha_1)^{16}\text{O}^*$ data, the α_1 channel may significantly contribute to the total astrophysical factor above about $E^{c.m.} = 50$ keV, calling for new data at low energies.

The most comprehensive R-matrix analysis, however not including the results in Zhang *et al.* (2021b), is the one carried out in deBoer *et al.* (2021b). For the α_1 channel, the results are consistent with Lombardo *et al.* (2019). For the $\alpha\gamma$ channel, the conclusions are that the α_2 channel is the major one and, as shown by Zhang *et al.* (2021b), low- and sub-threshold resonances can enhance the astrophysical factor to dominate over the α_0 channel at astrophysical energies. Finally, the result of the R-matrix analysis for the α_0 channel supports a flat S-factor devoid of resonances, in contrast with Indelicato *et al.* (2017); La Cognata *et al.* (2011, 2015a); Lombardo *et al.* (2013, 2015), in agreement with the analysis of Herndl *et al.* (1991). However, this is based on the data from Lorenz-Wirzba (1978), which were excluded from the NACRE compilation due to a likely large underestimate of reported absolute cross sections, by about a factor of 2.

In summary, while the situation for the α_0 and α_1 channels is well constrained, the contribution of higher ^{16}O excited states and, in particular, of the α_2 channel, needs further studies to confirm the occurrence of the $E_r^{c.m.} = 11$ keV resonance. This represents, at present, the largest source of uncertainty. Therefore, the recommended reaction rate lower limit is conservatively set by the Indelicato *et al.* (2017) one, while the upper limit is given by Zhang *et al.* (2021a), assuming the existence of the $E_r^{c.m.} = 11$ keV resonance.

D. Reactions of the neon-sodium (NeNa) cycle

1. $^{20}\text{Ne}(p, \gamma)^{21}\text{Na}$

The low energy S-factor of the $^{20}\text{Ne}(p, \gamma)^{21}\text{Na}$ reaction is dominated by the high energy tail of a subthreshold resonance only a few keV below the reaction threshold. In addition, the non-resonant capture into two additional states is relevant and provides the largest contribution in the energy range where data are available. All three transitions have been measured over a wide range of energies by two direct experiments (Lyons *et al.*, 2018; Rolfs *et al.*,

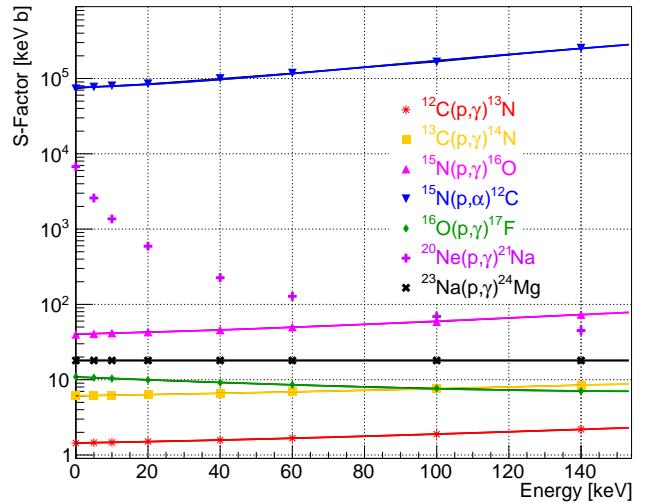


FIG. 10 Astrophysical S-factors of CNO and NeNa cycle reactions. The circles represent the values listed in Table XII. The solid lines represent polynomial (quadratic) fits of the S-factor values. The fit coefficients are listed in Table I. The S-factor of the $^{20}\text{Ne}(p, \gamma)^{21}\text{Na}$ reaction cannot be accurately fit with a quadratic (or cubic) polynomial function.

1975), and the results can be used to arrive at S-factor recommendations. However, no consistent parametrizations have been published that could be used to compare the results.

Here, the S-factor data from these two experiments were either extracted from the published supplemental information, or digitized from published figures and normalized to the strength of the $E_r^{c.m.} = 1113$ keV resonance ($\omega\gamma = 0.94$ eV; Lyons *et al.*, 2018). The resonant capture into the tail of the subthreshold resonance ($E_r^{c.m.} = -7.9$ keV; Firestone 2015) can be described by the Breit-Wigner equation, as shown in Iladis (2015); Rolfs and Rodney (1988). To account, additionally, for the non-resonant capture to the ground state, a constant value of $S(E) = 0.0463$ keVb, as recommended by the indirect experiment of Mukhamedzhanov *et al.* (2006), was included in the fit to the data of Lyons *et al.* (2018); Rolfs *et al.* (1975). For the two direct-capture transitions, the energy dependence of the S-factor was calculated with the program DIRCAP (Iladis, 2022), using a Woods-Saxon potential with a diffuseness of $a = 0.65$ fm to achieve S-factor magnitudes comparable to the experimental data. These S-factors were fit individually to the experimental data and then summed to the total S-factor. The results are listed in Table XII. Uncertainties from the individual contributions were added in quadrature.

However, significant and unquantifiable uncertainties likely exist because the results of the two experiments differ systematically both in magnitude and energy dependence (Lyons *et al.*, 2018). This may be caused by experimental or data-analysis errors in either experiment, or in both. In addition, the experiments at low ener-

gies only measured in a small angular range and relied on sparse angular distribution measurements at higher energies to infer angular correlation effects. Also, the experiments did not reach sufficiently low energies to conclusively determine the potential parameters to be used in the DIRCAP calculation. Consequently, this will contribute to the S-factor uncertainty when the energy range of stellar burning, depending on the temperature, is located far below the experimental range.

A new measurement of the $^{20}\text{Ne}(p, \gamma)^{21}\text{Na}$ direct-capture cross section down to $E^{c.m.} = 250$ keV has recently been published by the LUNA collaboration (Masha *et al.*, 2023). The potential-model results agree with the new experimental data.

2. $^{21}\text{Ne}(p, \gamma)^{22}\text{Na}$

At temperatures below 0.1 GK, the reaction rate is dominated by three narrow resonances at 16.6, 94.6 and 126 keV. The 16.6 and 94.6 keV resonances have never been observed directly in the (p, γ) channel (Görres *et al.*, 1982). Upper limits on their strengths were derived from a $^{21}\text{Ne}(d,p)^{22}\text{Na}$ experiment (Iliadis *et al.*, 2001; Neogy *et al.*, 1972). Results are reported in Table XIII. The non-resonant capture S-factor at zero energy was estimated by Rolfs *et al.* (1975) using single-particle spectroscopic factors of the bound states in ^{22}Na . According to the authors, the result, $S(0) \approx 20$ keV b, is uncertain by a factor of three.

3. $^{22}\text{Ne}(p, \gamma)^{23}\text{Na}$

The $^{22}\text{Ne}(p, \gamma)^{23}\text{Na}$ reaction used to be the most uncertain of the NeNa cycle, because of a number of poorly-known low-lying resonances. At temperatures below 30 MK, the reaction rate is dominated by a resonance at 36 keV. Its strength has been determined from a proton-transfer experiment (Hale *et al.*, 2001; Iliadis *et al.*, 2010). Two near-threshold resonances are expected at 5 and 28 keV, but the existing upper limits on their strengths are so low that they do not contribute significantly to the reaction rate (Hale *et al.*, 2001). Two supposed resonances at 68 and 100 keV, first reported in Powers *et al.* (1971), were discussed at length in the literature and searched for both with direct experiments (Cavanna *et al.*, 2015; Depalo *et al.*, 2016; Ferraro *et al.*, 2018), proton transfer reaction (Hale *et al.*, 2001) and inelastic scattering (Carrasco-Rojas *et al.*, 2023). None of such experiments lead to a positive detection of the two resonances. The latest direct upper limits on their strength suggest that, even if the two resonances exist, their contribution to the mean (recommended) thermonuclear reaction rate is negligible (Ferraro *et al.*, 2018). A resonance doublet dominates the reaction rate at temperatures around 0.1 GK.

The strength of the stronger resonance in the doublet was measured both with direct (Cavanna *et al.*, 2018; Kelly *et al.*, 2017; Lennarz *et al.*, 2020; Williams *et al.*, 2020) and proton-transfer experiments (data from Hale *et al.* (2001); Powers *et al.* (1971) re-analyzed in Santra *et al.* (2020)). The value adopted in Table XIII is an average of all literature values mentioned above. The strength of the weaker resonance in the doublet was inferred from proton-transfer data (Santra *et al.*, 2020).

At 0.2 GK, the reaction rate is mainly determined by a resonance at 181 keV. Its strength was measured in four independent experiments (Cavanna *et al.*, 2018; Ferraro *et al.*, 2018; Kelly *et al.*, 2017; Williams *et al.*, 2020). The strength recommended in Table XIII is a weighted average of all four values. The most intense higher-lying resonances, relevant to nova and supernova explosions, have been recently addressed in Depalo *et al.* (2015); Kelly *et al.* (2017); Longland *et al.* (2010); Williams *et al.* (2020). The non-resonant capture cross section has been measured by different groups (Ferraro *et al.*, 2018; Kelly *et al.*, 2017; Williams *et al.*, 2020). The low-energy trend of the ground state capture is described by a constant value of $S = (50 \pm 12)$ keVb, plus the contribution of a broad subthreshold state at -130 keV ($E_x = 8664$ keV) (Ferraro *et al.*, 2018).

4. $^{23}\text{Na}(p, \gamma)^{24}\text{Mg}$

The $^{23}\text{Na}(p, \gamma)^{24}\text{Mg}$ reaction has seen several recent studies (Boeltzig *et al.*, 2019, 2022; Caciolli *et al.*, 2011; Marshall *et al.*, 2021). The cross section is dominated by three narrow resonances over the lowest experimentally measured energy range, located at $E_r^{c.m.} = 133, 240.6,$ and 295.9 keV. Below the lowest known resonance, non-resonant components are thought to dominate.

Several indirect measurements have been made to characterize energies and proton spectroscopic factors of bound and unbound states in ^{24}Mg . The measurement of Hale *et al.* (2004) used the $^{23}\text{Na}(^3\text{He}, d)^{24}\text{Mg}$ reaction to investigate unbound states ($11690 < E_x < 12520$ keV) near the proton separation energy ($S_p = 11693$ keV). For lower-energy bound states that were not measured, the spectroscopic factors of Garrett *et al.* (1978) were adopted. A direct capture model was used to calculate the non-resonant component of the cross section. This was followed by an investigation of the lowest energy resonance at $E_r^{c.m.} \approx 133$ keV by Cesaratto *et al.* (2013), where an upper limit for the resonance strength was determined. A subsequent study at the LUNA facility (Boeltzig *et al.*, 2019) was able to make a first measurement of the resonance strength, $1.46_{-0.53}^{+0.58} \times 10^{-9}$ eV. A recent $^{23}\text{Na}(^3\text{He}, d)^{24}\text{Mg}$ transfer study at the Triangle Universities Nuclear Laboratory (TUNL) revised the energy of the ≈ 133 keV resonance to $E_r^{c.m.} = 133(3)$ keV (Marshall *et al.*, 2021).

Direct cross section measurements of the $^{23}\text{Na}(p, \gamma)^{24}\text{Mg}$ reaction have been made at the University of Notre Dame at energies between $E^{\text{lab}} = 500$ and 1000 keV by Boeltzig *et al.* (2022). Two, broad, strong, resonances were observed in several γ -ray transitions at $E_r^{\text{c.m.}} = 841$ and 981 keV whose low energy tails contribute in the low energy region. Note that the energy uncertainties presented here reflect the uncertainty in the beam calibration of the accelerator, which was neglected in Boeltzig *et al.* (2022). Further, measurements of the $E_x = 10740$ keV primary transition in ^{24}Mg , predicted to have the largest non-resonant capture component based on the spectroscopic factors of Garrett *et al.* (1978), observed a trend in the off-resonance cross section that was highly suggestive of non-resonant capture. Given the level of consistency observed between the non-resonant capture S-factors calculated from the spectroscopic factors of Garrett *et al.* (1978) and the external capture S-factors from the R-matrix analysis, an uncertainty in the non-resonant S-factor of 50% was recommended.

We have adopted the non-resonant component of Boeltzig *et al.* (2022) in this work, which, like all previous calculations, relies heavily on the spectroscopic factors of Garrett *et al.* (1978).

5. $^{23}\text{Na}(p, \alpha)^{20}\text{Ne}$

The $^{23}\text{Na}(p, \alpha)^{20}\text{Ne}$ reaction rate is dominated by narrow resonances, with a total of 52 resonances in the center-of-mass energy interval 6 – 2328 keV.

After NACRE, the state-of-the-art tabulation of the resonance strengths is the one in Iliadis *et al.* (2010). It is mostly based on the ($^3\text{He}, \text{d}$) proton transfer measurement by Hale *et al.* (2004), supplying both the excitation energies and spectroscopic factors of the near-threshold states. In detail, below $E^{\text{c.m.}} = 968$ keV, all the resonance strengths are taken from Table VI of Hale *et al.* (2004), with the exception of the $E_r^{\text{c.m.}} = 133$ keV resonance strength (Marshall *et al.*, 2021), for which the upper limit directly determined by Iliadis *et al.* (2005); Rowland *et al.* (2004) is adopted, equal to 1.5×10^{-8} eV. Above $E^{\text{c.m.}} = 968$ keV, the Hale *et al.* (2004) strengths are rescaled to the one of the $E_r^{\text{c.m.}} = 338$ keV resonance determined by Rowland *et al.* (2002). This result is considered as reference value since it provides the most recent directly-measured strength.

Besides the above-mentioned works, two more should be considered. One is Iliadis *et al.* (2005), which is not used in Iliadis *et al.* (2010) because the same upper limit for the $E_r^{\text{c.m.}} = 133$ keV resonance as in Rowland *et al.* (2004) is reported. The other work is Cesaratto *et al.* (2013), which used the upper limit on the (p, γ) resonance strength and the upper limit on the α branching ratio (equal to 0.13 at the 95% CL), to provide an up-

per limit on the (p, α) strength of the $E_r^{\text{c.m.}} = 133$ keV resonance equal to 0.88×10^{-9} eV. This implies a negligible contribution to the reaction rate and confirms the conclusions drawn in Iliadis *et al.* (2010).

XII. Electron screening of nuclear reactions

Nuclear fusion reactions measured in the laboratory and those occurring in the solar core are both affected by the electronic environments in which they take place. The effects are, however, different requiring special care in deriving the appropriate solar reaction rates. As long as these rates cannot be measured under solar plasma conditions in the laboratory, the usual two-step strategy is to remove the laboratory screening effects from the data to obtain the bare nuclear cross section σ^b , which is then modified to incorporate the solar plasma screening modifications. Thus, one must address the differential effects of screening for the solar cross section σ^{solar} and the laboratory results, σ^{lab} .

A. Screening in laboratory experiments

Screening diminishes the Coulomb barriers that retard interactions between bare nuclei, thus enhancing cross sections, with the effects becoming more pronounced with decreasing center-of-mass energy. Nuclear astrophysics has advanced rapidly over the last two decades with the deployment of underground accelerators and other low-background techniques (Broggini *et al.*, 2010) that have enabled low-energy, low-counting-rate experiments. In particular, laboratory measurements of the pp-chain reactions $^2\text{H}(p, \gamma)^3\text{He}$ and $^3\text{He}(^3\text{He}, 2p)^4\text{He}$ have been made at energies corresponding to the solar Gamow peak. A treatment of the distorting effects of screening then becomes quite important in extracting the bare S-factor $S^b(E)$.

The bare cross section is given by

$$\sigma^b(E) = \frac{1}{E} S(E) P(E); \quad P(E) \equiv \exp(-2\pi\eta(E)) \quad (24)$$

with $\eta(E) = Z_1 Z_2 \alpha \sqrt{\mu/2E}$, where μ is the reduced mass and E the center-of-mass energy. The effects of screening can be incorporated into the Gamow penetration factor $P(E)$ through an energy shift, $E \rightarrow E + U_e$, where the screening potential U_e has the net effect of lowering the Coulomb barrier. At the lowest accessible energies in laboratory measurements, the separation of target and projectile during tunnelling is much smaller than the atomic radius: thus the electrons, originally bound to the target (assuming a projectile beam of bare nuclei), are attracted by the joint charge of target and projectile during the fusion. The screening potential can then be replaced by a constant screening energy U_e . This picture was confirmed

TABLE XI Overview of CNO and NeNa reactions.

Reaction	Category ^a	$S(0) \pm \Delta S(0)$ (keVb)	E_0 ^b (keV)	Most recent reference	Comments
$^{12}\text{C}(p,\gamma)^{13}\text{N}$	B	1.44 ± 0.06	25	Skowronski <i>et al.</i> (2023a)	SF II: $S(0) = 1.34 \pm 0.21$ keVb
$^{13}\text{C}(p,\gamma)^{14}\text{N}$	B	6.1 ± 0.4	25	Skowronski <i>et al.</i> (2023a)	SF II: $S(0) = 7.6 \pm 1.0$ keVb
$^{15}\text{N}(p,\gamma)^{16}\text{O}$	B	40 ± 3	28	deBoer <i>et al.</i> (2013)	SF II: $S(0) = 36 \pm 6$ keVb
$^{15}\text{N}(p,\alpha)^{12}\text{C}$	B	73000 ± 5000	28	La Cognata <i>et al.</i> (2009)	SF II: $S(0) = 73000 \pm 5000$ keVb
$^{16}\text{O}(p,\gamma)^{17}\text{F}$	B	10.92 ± 0.44	30	Iliadis <i>et al.</i> (2022)	SF II: $S(0) = 10.6 \pm 0.8$ keVb
$^{17}\text{O}(p,\gamma)^{18}\text{F}$	B	4.7 ± 1.0	31	Gyürky <i>et al.</i> (2017)	SF II: $S(0) = 6.2 \pm 3.1$ keVb
$^{17}\text{O}(p,\alpha)^{14}\text{N}$	A	—	31	Bruno <i>et al.</i> (2016)	near-threshold resonances
$^{18}\text{O}(p,\gamma)^{19}\text{F}$	B	23.0 ± 3.8	31	Pantaleo <i>et al.</i> (2021)	also in literature: $S(0) = 7.1$ and 15.7 keVb
$^{18}\text{O}(p,\alpha)^{15}\text{N}$	B	—	31	Bruno <i>et al.</i> (2019)	low-energy resonances
$^{19}\text{F}(p,\gamma)^{20}\text{Ne}$	A	—	33	Zhang <i>et al.</i> (2022)	possible resonance near threshold
$^{19}\text{F}(p,\alpha)^{16}\text{O}$	B	—	33	Zhang <i>et al.</i> (2021a)	possible resonance near threshold
$^{20}\text{Ne}(p,\gamma)^{21}\text{Na}$	B	6776 ± 550	36	present work	
$^{21}\text{Ne}(p,\gamma)^{22}\text{Na}$	A	≈ 20	36	Rolfs <i>et al.</i> (1975)	factor 3 uncertainty in $S(0)$
$^{22}\text{Ne}(p,\gamma)^{23}\text{Na}$	A	415 ± 91	36	Ferraro <i>et al.</i> (2018)	sub-threshold resonance at -130 keV
$^{23}\text{Na}(p,\gamma)^{24}\text{Mg}$	B	18 ± 9	38	Boeltzig <i>et al.</i> (2022)	
$^{23}\text{Na}(p,\alpha)^{20}\text{Ne}$	A	—	38	Iliadis <i>et al.</i> (2010)	

^a Categories for the dominant rate contribution at the central temperature of the Sun: (“A”) Narrow resonances. (“B”) Broad resonant/non-resonant contributions. We define a “narrow resonance” by the criterion that its full contribution to the reaction rate can be calculated from the resonance energy and strength alone.

^b Central energy location (center-of-mass system) of Gamow peak at the solar core.

TABLE XII S-factors versus center-of-mass energy. Data are also plotted in Fig. 10.

Reaction	$S(E)$ (keVb) (% uncertainty)						
	5 keV	10 keV	20 keV	40 keV	60 keV	100 keV	140 keV
$^{12}\text{C}(p,\gamma)^{13}\text{N}$ ^a	$1.46(4.1\%)$	$1.47(4.1\%)$	$1.51(4.1\%)$	$1.58(4.1\%)$	$1.67(4.1\%)$	$1.89(4.0\%)$	$2.19(4.1\%)$
$^{13}\text{C}(p,\gamma)^{14}\text{N}$ ^a	$6.16(6.0\%)$	$6.22(6.0\%)$	$6.34(6.1\%)$	$6.60(6.1\%)$	$6.89(6.1\%)$	$7.58(6.1\%)$	$8.47(6.1\%)$
$^{15}\text{N}(p,\gamma)^{16}\text{O}$	$40.8(7.5\%)$	$41.7(7.5\%)$	$43.0(7.5\%)$	$45.8(7.5\%)$	$50.0(7.5\%)$	$59.0(7.5\%)$	$73.4(7.5\%)$
$^{15}\text{N}(p,\alpha)^{12}\text{C}$	$7.7 \times 10^4(6.8\%)$	$8.0 \times 10^4(6.8\%)$	$8.5 \times 10^4(6.8\%)$	$1.00 \times 10^5(6.8\%)$	$1.18 \times 10^5(6.8\%)$	$1.66 \times 10^5(6.8\%)$	$2.53 \times 10^5(6.8\%)$
$^{16}\text{O}(p,\gamma)^{17}\text{F}$ ^b	$10.64(4.0\%)$	$10.38(4.0\%)$	$9.92(4.0\%)$	$9.15(4.0\%)$	$8.56(4.0\%)$	$7.67(4.0\%)$	$7.05(4.0\%)$
$^{20}\text{Ne}(p,\gamma)^{21}\text{Na}$ ^c	$2586(8\%)$	$1365(8\%)$	$593(7\%)$	$226(7\%)$	$128(7\%)$	$69(6\%)$	$45(6\%)$
$^{23}\text{Na}(p,\gamma)^{24}\text{Mg}$	$18(50\%)$	$18(50\%)$	$18(50\%)$	$18(50\%)$	$18(50\%)$	$18(50\%)$	$18(50\%)$

^a From Skowronski *et al.* (2023a).

^b From Iliadis *et al.* (2022).

^c From present work.

in simulations of the electron dynamics in low-energy fusion reactions (Shoppa *et al.*, 1993). After expanding the argument of the exponential in Eq. 24 to first order in U_e one obtains the following relation between the laboratory and bare S-factors (Assenbaum *et al.*, 1987)

$$S^{\text{lab}}(E) = S^b(E) \exp\left(\frac{\pi\eta(E)U_e}{E}\right) \quad (25)$$

This result neglects a small normalization correction associated with the effects U_e on the matching of the external wave function to that in the Coulomb region.

The $1/E$ dependence of the exponent is responsible for the growing importance of the screening correction at low energies. Although the screening potential U_e in principle

can be calculated from the electron charge distribution, in most applications it is treated as a free constant parameter, determined along with $S^b(0)$ and its derivatives from a fit to data. Figure 4 of SF II shows the results for $S_{33}(E)$, where the fit yielded $U_e = 305 \pm 90$ eV. The exponential dependence of the screening correction leads to a rapid increase in $S_{33}^{\text{lab}}(E)$ relative to S_{33}^b , reaching $\sim 40\%$, at the lowest data point measured. This example also underscores the potential for screening to complicate the extraction of $S^b(E)$ for energies relevant to the solar core, even in cases where measurements are restricted to higher energy than the Gamow window.

In principle one can test the adequacy of Eq. 25 by doing laboratory experiments at low energies – even if

TABLE XIII Recommended strengths and relative kinetic energies of narrow resonances.

Reaction	$E_{\text{c.m.}}^{\text{c.m.}}$ (keV)	$\omega\gamma$ (eV) ^b	Comments/references
$^{17}\text{O}(p,\gamma)^{18}\text{F}$	65	$15.6^{+18.3}_{-4.0} \times 10^{-12}$	see text for details
$^{17}\text{O}(p,\alpha)^{14}\text{N}$	183	$(1.77 \pm 0.09) \times 10^{-6}$	see text for details
	65	$4.6^{+5.4}_{-1.2} \times 10^{-9}$	see text for details
$^{18}\text{O}(p,\gamma)^{19}\text{F}$	183.9	$(1.68 \pm 0.03_{\text{stat}} \pm 0.12_{\text{syst}}) \times 10^{-3}$	Bruno <i>et al.</i> (2015); Chafa <i>et al.</i> (2007)
	20	5.7×10^{-22}	Champagne and Pitt (1986)
	90	$(0.53 \pm 0.07_{\text{stat}} \pm 0.07_{\text{syst}}) \times 10^{-9}$	Best <i>et al.</i> (2019)
	143	$(0.98 \pm 0.03) \times 10^{-3}$	see text for details
$^{18}\text{O}(p,\alpha)^{15}\text{N}$	20	$8.3^{+3.8}_{-2.6} \times 10^{-19}$	La Cognata <i>et al.</i> (2008)
	90	$2.0^{+13.7}_{-0.4} \times 10^{-7}$	see text for details
	143	$(166 \pm 8) \times 10^{-3}$	see text for details
$^{19}\text{F}(p,\gamma)^{20}\text{Ne}$	11.0	$\leq 8.5 \times 10^{-29}$	Betts <i>et al.</i> (1975); Kious (1990)
	212.7	$\leq 4.2 \times 10^{-6}$	Spyrou <i>et al.</i> (2000); Zhang <i>et al.</i> (2022)
	225.2	$(4.19 \pm 0.33) \times 10^{-5}$	Spyrou <i>et al.</i> (2000); Zhang <i>et al.</i> (2022)
	323.9	$(3.16 \pm 0.33) \times 10^{-3}$	Couture <i>et al.</i> (2008); Williams <i>et al.</i> (2021); Zhang <i>et al.</i> (2022)
$^{19}\text{F}(p,\alpha\gamma)^{16}\text{O}$	11.0	$(7.5 \pm 3.0) \times 10^{-29}$	Zhang <i>et al.</i> (2021a)
	212.7	0.0126 ± 0.0013	Spyrou <i>et al.</i> (2000); Zhang <i>et al.</i> (2021a)
	225.2	0.0011 ± 0.0004	Spyrou <i>et al.</i> (2000); Zhang <i>et al.</i> (2021a)
	323.9	23.5 ± 0.6	Zhang <i>et al.</i> (2022, 2021a)
$^{21}\text{Ne}(p,\gamma)^{22}\text{Na}$	16.6	$\leq 6.2 \times 10^{-24}$	Iliadis <i>et al.</i> (2001)
	94.6	$\leq 6.4 \times 10^{-10}$	Iliadis <i>et al.</i> (2001)
	126	$(3.75 \pm 0.75) \times 10^{-5}$	Görres <i>et al.</i> (1982)
$^{22}\text{Ne}(p,\gamma)^{23}\text{Na}$	5	$\leq 2.1 \times 10^{-51}$	Hale <i>et al.</i> (2001)
	28	$\leq 3.2 \times 10^{-25}$	Hale <i>et al.</i> (2001)
	35	$(3.1 \pm 1.2) \times 10^{-15}$	Iliadis <i>et al.</i> (2010)
	68	$\leq 6 \times 10^{-11}$	Ferraro <i>et al.</i> (2018)
	100	$\leq 7 \times 10^{-11}$	Ferraro <i>et al.</i> (2018)
	150	$(3.9 \pm 0.9) \times 10^{-9}$	Santra <i>et al.</i> (2020)
	151	$(1.97 \pm 0.12) \times 10^{-7}$	see text for details
	181	$(2.39 \pm 0.12) \times 10^{-6}$	see text for details
$^{23}\text{Na}(p,\gamma)^{24}\text{Mg}$	133	$1.46^{+0.58}_{-0.53} \times 10^{-9}$	Boeltzig <i>et al.</i> (2019); Marshall <i>et al.</i> (2021)
	240.6	$5.3(1) \times 10^{-4}$	Keinonen <i>et al.</i> (1989); Switkowski <i>et al.</i> (1975)
	295.90	$1.05(19) \times 10^{-1}$	Endt <i>et al.</i> (1990); Switkowski <i>et al.</i> (1975)
$^{23}\text{Na}(p,\alpha)^{20}\text{Ne}$	133	$\leq 1.5 \times 10^{-8}$	Iliadis <i>et al.</i> (2005); Marshall <i>et al.</i> (2021); Rowland <i>et al.</i> (2004)
	240.6	≤ 0.1	Hale <i>et al.</i> (2001)
	295.90	$1.03(26) \times 10^{-2}$	Iliadis <i>et al.</i> (2010)

^a We define a “narrow resonance” by the criterion that its full contribution to the total rate can be calculated from the resonance energy and strength alone.

^b The resonance strength, in the center-of-mass system, for the reactions listed here is defined as $\omega\gamma \equiv (2J+1)(2j_t+1)^{-1}(2j_p+1)^{-1}\Gamma_p\Gamma_x/\Gamma$, with J , j_t , and j_p the spins of the resonance, target, and projectile, respectively, and Γ_p , Γ_x , and Γ the proton partial width, γ -ray or α -particle partial width, and total resonance width, respectively.

the reactions studied are not directly relevant to nuclear astrophysics. Such studies began more than 30 years ago (Assenbaum *et al.*, 1987), using gaseous targets to limit energy loss effects and other systematics. Subsequent studies have varied greatly in their configurations, e.g., targets ranging from atomic and molecular gases to metals,¹⁰ thin and thick targets, and use of direct and inverse kinematics (Aliotta *et al.*, 2022b). The experiments can be very challenging at the energies needed, subject to

backgrounds from both intrinsic activities in detectors and external sources such as cosmic rays.

The values of U_e extracted from experiment using gaseous targets have often exceeded the adiabatic limit, defined in atomic physics as the difference between the electron binding energies of the separate atoms in the entrance channel and that of the composite atom (Aliotta *et al.*, 2001; Angulo *et al.*, 1993; Engstler *et al.*, 1988, 1992; Greife *et al.*, 1995; Prati *et al.*, 1994; Rolfs, 2001; Rolfs and Somorjai, 1995). This in turn has generated unease about the reliability of the values for $S^b(E)$ extracted from laboratory measurements for use in astrophysics (Balantekin *et al.*, 1997).

¹⁰ In this section, the word “metals” is used with the traditional meaning from chemistry, not the astrophysical one.

Several suggestions have been made to resolve this apparent discrepancy. For example, an error in the estimated stopping power could result in a smaller reaction effective energy, so that the experimentally deduced screening would be larger than the true value (Bang *et al.*, 1996; Langanke *et al.*, 1996). Indeed such deviations have been found between theoretical calculations of stopping powers and the tabulation which is traditionally being used in the analysis of low-energy fusion data (Bertulani, 2004; Bertulani and de Paula, 2000; Golser and Semrad, 1991). An accurate experimental determination of stopping powers for gaseous hydrogen and helium targets would be very desirable. Corrections associated with the non-uniform distribution of the electron cloud around the nucleus (Carraro *et al.*, 1988) and the effects of electron-electron interactions (Shoppa *et al.*, 1993) have also been suggested as potential generators of unidentified systematic error. For molecular targets the screening energy shows a significant dependence on the scattering angle, being smallest if the projectile passes the spectator nucleus before fusion (Shoppa *et al.*, 1996). Furthermore, the experimental deduction of screening energies requires assumptions about the bare nuclear cross sections, which is another source of uncertainty. This was demonstrated in Tumino *et al.* (2021, 2014), where after constraining fits by the addition of higher energy data (from a Trojan Horse measurement), a lower value of U_e was obtained. Clusterization and polarization effects have also been proposed to effect the deduced screening energies in specific low-energy reactions (Spitaleri *et al.*, 2016, 2017)

For resonances, both the energy and width can be affected by screening (Salpeter, 1954; Zinner, 2007). Particular care has to be taken for narrow, low-energy resonances where the assumption of a constant screening energy is not valid and the radial dependence of the screening potential has to be considered which leads to a significant reduction of screening on the resonance width (Iliadis, 2023).

Particularly large screening energies were observed in deuteron- and proton-induced fusion reactions if the projectile was implanted into metals (Czerski *et al.*, 2001; Raiola *et al.*, 2002a,b; Kasagi *et al.*, 2002; Czerski *et al.*, 2006; Cvetinovic *et al.*, 2015; Lipoglavsek and Cvetinović, 2020). While the d+d fusion data obtained with a gaseous target (Greife *et al.*, 1995) were compatible with the adiabatic limit ($U_e = 20$ eV, Bracci *et al.*, 1991, 1990), screening energies obtained for various metal hosts varied among themselves and could reach values of several 100 eV. As the adiabatic limit derived for the fusion on individual atoms in a gas, does not apply to reactions with targets implanted in a host material, an appropriate model has been developed (Czerski *et al.*, 2004; Huke *et al.*, 2008) which agrees nicely with recent screening energies obtained under high-vacuum conditions (Czerski, 2022; Czerski *et al.*, 2016). Huke *et al.* (2008) also identified oxygen and carbon contamination as the source for

some particularly large deduced screening energies. We note that the studies of d+d low-energy fusion were motivated by the quest to enhance nuclear fusion rates by environmental effects (Czerski, 2022).

The potential discrepancy between observed and theoretical screening energies would become astrophysically irrelevant if it were possible to measure reaction rates directly under solar plasma conditions. This may prove possible using high-intensity lasers. First measurements of light-ion reactions in plasma have been performed, though under plasma conditions for which screening effects are expected to be negligible (Barbui *et al.*, 2013; Casey *et al.*, 2023; Zylstra *et al.*, 2016, 2020). A direct measurement of bare nuclear cross section, thus avoiding screening effects, is proposed for storage ring experiments, in which a stored beam of ions can collide with a transverse beam (Glorius and Bruno, 2023).

B. Screening in the solar core

The energy-generating reactions that occur in the solar core involve nuclei that are almost completely ionized. Consequently the atomic environment differs substantially from that of the terrestrial experiments in which these same reactions are measured. Once $S^b(E)$ is determined from these laboratory measurements, the screening effects of the plasma must be folded in to yield $S^{\text{solar}}(e)$. Adopting the notation of SF II, the correction factor is defined as

$$f_0(E) = \frac{\sigma^{\text{solar}}(E)}{\sigma^b(E)} \quad (26)$$

In the weak screening approximation (Salpeter, 1954), the ion-ion Coulomb potential is screened on a length scale given by the Debye radius R_D

$$V(r) = \frac{\alpha Z_1 Z_2}{r} \exp\left(-\frac{r}{R_D}\right) \quad (27)$$

yielding

$$f_0 \sim \exp\left(\frac{\alpha Z_1 Z_2}{R_D kT}\right) \quad (28)$$

We refer the reader to SF II for the functional dependence of R_D on temperature, density, and the mass fractions X_i , and for discussions of corrections for electron degeneracy and incomplete ionization.

SF II discusses the conditions under which the weak screening approximation is valid: under solar core conditions this requires $Z_1 Z_2 < 10$, a condition satisfied by pp-chain and CNO bi-cycle reactions. Yet corrections are expected at some level, as the inter-ion potential is not completely negligible, in all collisions, compared to the relative kinetic energy of the ions (Bahcall *et al.*, 2002; Brown and Sawyer, 1997; Carraro *et al.*, 1988; Fiorentini

et al., 2003; Itoh *et al.*, 1979; Ogata, 1997; Salpeter, 1954; Salpeter and van Horn, 1969).

Dynamic corrections –nonadiabatic effects that arise when the velocities of reacting nuclei momentarily exceed typical plasma velocities– have been a point of some controversy that both SF I and SF II addressed. However, the absence of such dynamic corrections even for large Gamow energies has been shown to be a consequence of the nearly exact thermodynamic equilibrium of the solar plasma (Brown and Sawyer, 1997; Gruzinov, 1998; Gruzinov and Bahcall, 1998). Specifically, Bahcall *et al.* (2002) found that corrections to the Salpeter formula under solar conditions would be at the \sim few per cent level, and pointed to specific errors in several papers that had come to contrary conclusions.

There have been a few recent papers advocating changes in solar rates due to dynamical corrections, including Vescovi *et al.* (2019), Wood *et al.* (2018), Simonucci *et al.* (2013), Mussack (2011), Mussack and Däppen (2010), and Mao *et al.* (2009). The approaches employed are based on modeling the plasma, and in general the authors do not relate their conclusions to earlier work, particularly Bahcall *et al.* (2002), making them difficult to evaluate. The essential point of Bahcall *et al.* (2002) is the conceptual simplifications that result from recognizing that screening in the solar plasma can be formulated as a problem in equilibrium statistical mechanics. This approach removes all need to classified rates as fast or slow relevant to some plasma timescale, as there is no time in equilibrium statistical mechanics. Consequently we continue to regard Bahcall *et al.* (2002) as the most realistic estimate of corrections to the Salpeter formula.

XIII. Radiative opacities

A. Introduction

Opacity is a measure of the photon absorption of matter and is an essential quantity for understanding radiative heat transfer in the Sun. Radiative heat transfer occurs through the absorption and emission processes taking place within the material that the radiation traverses. In local thermodynamic equilibrium (LTE)¹¹, the emission and absorption are straightforwardly related through

¹¹ Local thermodynamic equilibrium can be defined as the situation where the thermodynamic properties of a microscopic volume of matter are the same as their thermodynamic equilibrium values corresponding to the local electron temperature and density. The assumption of LTE is commonly used in stellar modeling, and the discussion of opacities in this section will be limited only to LTE conditions. In some circumstances (such as solar coronal modeling), non-LTE conditions should be given careful consideration, but this will not be discussed in this paper.

the Planck function. In this scenario, determining the absorption coefficients or opacity, denoted as κ_ν , provides a complete description of how radiation is transported through the material.

For example, in the equilibrium diffusion limit, the radiation heat flux F_R is directly related to the Rosseland-mean opacity (RMO) κ_R through the following equation (Huebner and Barfield, 2014)

$$F_R = -\frac{16\pi}{3} \frac{\sigma T^3}{\kappa_R} \nabla T, \quad (29)$$

where T represents temperature, ∇T denotes its gradient, and σ is the Stefan-Boltzmann constant. The quantity κ_R is the solar Rosseland mean opacity (RMO). Thus, any error in this quantity introduces error in the radiative heat transport and the simulated solar evolution. Quantifying the error in κ_R is very challenging because κ_R is a weighted mean of a complex solar mixture opacity:

$$\frac{1}{\kappa_R} = \int_0^\infty \frac{1}{\kappa_\nu} w_\nu d\nu, \quad (30)$$

where w_ν is a weighting function related to the temperature derivative of the Planck function B_ν :

$$B_\nu = \frac{2h\nu^3}{c^2} \left(e^{h\nu/kT} \right)^{-1}, \quad (31)$$

where ν denotes the photon frequency, h the Planck constant, k the Boltzmann constant, and c the speed of light. Moreover, the frequency-resolved solar opacity κ_ν is a sum of elemental opacities $\kappa_{\nu,i}$ weighted by their mass fractions (b_i):

$$\kappa_\nu = \sum_i b_i \kappa_{\nu,i} \quad (32)$$

Equation 30 reveals that the accuracy of κ_ν is especially important over the spectral range where w_ν is high and κ_ν is low (due to the inverse weighting). Equation 32 also indicates that accurate κ_ν is possible only when we have both accurate element opacities $\kappa_{\nu,i}$ and accurate abundances b_i .

Figure 11 shows the total monochromatic opacity of the Sun at three different solar radii, highlighting the variations in w_ν (gray), κ_ν (black), and some of the leading individual element contributions $\kappa_{\nu,i}$ (red, green, blue) to the RMO determination (Eq. 30), as a function of radius. This figure was made assuming the GS98 solar abundances and using OPLIB opacities (Colgan *et al.*, 2016), and is included primarily for illustrative purposes. As shown in the figure, the solar opacity varies considerably in both magnitude (note the log axes) and shape across different solar radii. Moreover, each element contributes distinctively to the overall monochromatic opacity, with either a dominant continuous opacity spectrum,

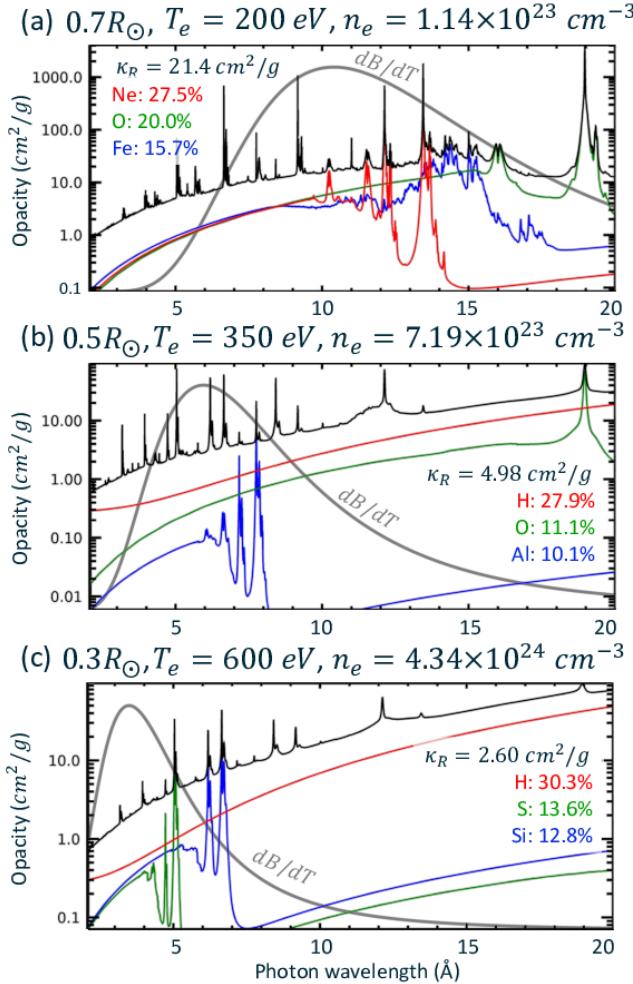


FIG. 11 Monochromatic opacity of the Sun at three different solar radii as indicated. The temperature and electron density at each solar radius are displayed, as well as some of the major elemental opacity contributions at these conditions. Note that not all contributions are plotted for clarity. The percentages of three elements displayed indicate the leading elemental contributors to the Rosseland mean opacity (RMO) at these conditions. These are roughly quantified by calculating how much the RMO is reduced if each element is completely removed. This figure was made using the GS98 solar abundances and OPLIB opacities, and is mainly intended to show the spectral complexity of the solar opacity.

or strong lines that arise from bound-bound transitions. If there is an error in the calculated solar opacity κ_ν , it would stem from inaccuracies in the calculated element opacities and/or abundances. Additionally, the error is expected to be a complex function of photon energy and would vary at each radius due to changes in temperature and density. As a result, it is not possible to provide a single factor for uncertainty or correction. It is crucial to assess the accuracy of the intricate spectral features depicted in Fig. 11 in order to better understand the nature of the errors.

These complex spectral features arise from microscopic

physics at the atomic level – that is, the absorption of radiation through quantum mechanical processes. If one knows the atomic energy levels and transition probabilities of all the relevant atomic states (of each atom or atomic ion of each material in the Sun), then the opacity of that material may be computed.¹² Since atomic structure and transition probabilities can be predicted (under various approximations), starting from the Schrödinger (or Dirac) equation, this seems like an achievable goal. However, complications quickly arise from a multitude of factors, such as: how to deal with the infinity of atomic states that may be populated in each atom or atomic ion; how to account for the perturbations caused by the plasma environment on the atomic states; and how to handle the large quantities of data that arise with such calculations, particularly for multi-electron atoms and ions. In solar modeling, one requires the opacity of all relevant elements (around 30) at all (plasma) conditions that are found at all radii throughout the solar evolution. Thus, the opacity must be tabulated as a function of species, plasma temperature, plasma density, and photon energy.

Nevertheless, steady progress has been made in determining the opacity of solar mixtures over the last five decades. In the 1980s, insufficient accuracy in stellar-envelope opacity was hypothesized through multiple astrophysical puzzles (Simon, 1982), and opacity models were subsequently improved by incorporating more detailed atomic physics and a more complete set of states (Berrington *et al.*, 1987; Iglesias and Rogers, 1996; Magee *et al.*, 1995; Seaton, 1987). The improved accuracy was confirmed through a few benchmark experiments (Davidson *et al.*, 1988; Perry *et al.*, 1991; Springer *et al.*, 1992; Winhart *et al.*, 1996) and improved agreement with some astrophysical observations (Daszyńska-Daszkiewicz and Walczak, 2009, 2010; Iglesias and Rogers, 1991). However, the calculated solar opacities are not experimentally validated at most radii, and in particular more deeply into the Sun where the conditions (temperature and density) are more extreme.

In the last decade, experimental methods have steadily improved, in concert with the availability of powerful x-ray sources, such as the Sandia Z facility. Monochromatic opacities were successfully measured at multiple conditions relevant to solar convection-base conditions (Bailey *et al.*, 2015, 2007). Systematic experimental studies performed on chromium, iron, and nickel helped narrow down hypotheses for the discrepancies (Nagayama *et al.*, 2019). All of these experimental results provided essential clues for testing various approximations for so-

¹² We note that the opacity of molecules and dust, or grains, can be very important in some circumstances (Huebner and Barfield, 2014); however they play a minor role in stellar modeling and so will not be discussed here.

lar interior opacity calculations; however, the reported model-data disagreements also raised significant controversy. This led to various theoretical investigations and the development of independent experimental methods (Opachich *et al.*, 2022; Perry *et al.*, 2020). Ongoing theoretical and experimental investigations are crucial for accurately interpreting the reported disagreements and to reach a consensus in this field.

In this section, our aim is not to provide a consensus but rather to summarize the history and challenges associated with solar opacity calculations and experiments, as well as provide an overview of the current status of stellar opacity investigations.

B. Opacity models

Over the years, a wide range of approximations has been used for the various components of opacity calculations. To help navigate these approximations, it may be useful to briefly summarize the key components of a frequency-dependent opacity. Opacities are often broken down into various components. The first of these is “bound-bound” absorption (a photoexcitation transition between two levels of an atomic ion), which manifests as strong line features in the opacity spectrum (such as the black ‘spikes’ in Fig. 11). The “bound-free” absorption (photoionization of the ground or excited level of an atomic ion) is normally a continuous (wavelength-dependent) contribution (such as the smooth (red) hydrogen opacity contribution in the middle and lower panels of Fig. 11). The “free-free” absorption (also known as inverse Bremsstrahlung) is where a free electron absorbs a photon, and finally “scattering” (the scattering of a photon by a free electron) contributions can be important at high photon energies. Other contributions may be important in molecules or other material. Equation 30 shows that due to the harmonic mean inherent in the RMO, the most important photon energy range is when the spectral opacity is low and when the weighting function w_ν is high. In the solar case, at a given solar radius, it is important to know which elements dominate the opacity at such a region, which will of course depend on the elemental abundance and its contribution at that distance, as illustrated in Fig. 11.

The determination of opacities usually starts from one of two fundamental approaches. The first approach, which was used in the first opacity computations (Huebner and Barfield, 2014), starts from a “mean ion” model, where a (fictitious) “average” atom or ion of a given element is conceived, with fractional occupation numbers, which is consistent with the given temperature and density of the species under consideration. The fractional occupation numbers, usually determined through a self-consistent iterative approach, inform the real (physical) ion stages of importance at such conditions. The average

atom can then be unfolded through various procedures to the physical ion stages of relevance, and the opacity of such ions can be determined. In this model the plasma effects (in particular the perturbation of the average atom by the fields produced by the surrounding plasma ions and electrons) can be included in a natural way, and this approach is often most useful at high densities, where plasma perturbations on the atoms or ions are most important. The infinity of atomic states are naturally truncated by plasma effects that remove bound states from consideration. The old Inferno model (Liberman, 1979) and the more recent STAR tables (Krief *et al.*, 2016b, 2018a) are based on this method. We note that very recent work (Gill *et al.*, 2023) has considerably improved some of the approximations used in related approaches.

The second approach is often referred to as detailed configuration (or term) accounting, where one determines all possible ion stages of the species and then determines all possible ground and (multiply) excited states of each ion, each with a corresponding population that is condition dependent. This method is often more spectroscopically accurate than the mean ion model, because it considers the structure of each ion separately, but can require enormous computational resources for complex species. The structure of each ion is determined from solution of the Schrödinger (or Dirac) equation, which can then also produce transition probabilities. Within this procedure, many computational difficulties must be overcome to produce acceptably accurate data for all ion stages of all relevant elements. A few examples are given for illustration. Within the solution of the Schrödinger equation it is desirable to include “configuration-interaction” (CI), a two-body term in the Hamiltonian that accounts for electron-electron interactions (Cowan, 1981), which results in a quite accurate description of atomic energy levels. However, the inclusion of this term for atomic ions with large numbers of configurations can quickly result in a computationally intractable problem. Therefore, approximations in which CI is included in only a limited manner are often employed. Furthermore, careful consideration has to be given to the number of states that should be included in the calculations. A compromise is always necessary between including as many states as possible versus the computational resources available for the calculation. Convergence checks are necessary to ensure that sufficient states are included. However, the rate of convergence of the opacity with respect to the number of states included will depend on the plasma conditions of interest, so this in itself is not necessarily a straightforward matter. The computer codes that are used to generate the opacity must also be able to efficiently process large amounts of atomic data.

Similarly, when considering photoexcitation and photoionization of atomic ions, it is desirable to include coupling of the bound electrons to the continuum electron(s). Often termed “close-coupling”, this may be ac-

complished through techniques such as the R-matrix approach (Burke, 2011; Seaton, 1987) or other related methods. However, such approaches are also computationally intensive, and can be severely limited in the number of terms that can be included in the resulting close-coupling expansions. It is also not clear how to include plasma perturbations within such approaches. As a result, perturbative approaches, such as distorted-wave methods (Badnell *et al.*, 2005; Cowan, 1981; Sampson *et al.*, 2009), are often employed, which are much more computationally tractable, and of sufficient accuracy for mid- and highly-ionized ions.

In concert with an atomic model, the thermodynamics of the system must be considered. That is, an equation-of-state (EOS) description of the material is required so that the thermodynamic properties and atomic state populations can be determined for a given material temperature and density. In the mean ion approach the thermodynamic quantities arise naturally through the description of the atom/ion in a plasma. In a Detailed Configuration Accounting approach, a common approach to EOS calculations starts from a chemical picture based on a minimization of the chemical free energy (Däppen *et al.*, 1987; Hakel and Kilcrease, 2004; Hummer and Mihalas, 1988; Rogers and Iglesias, 1992). Such a model has to treat a wide range of physical conditions, ranging from ideal gas through high density conditions where pressure ionization may rapidly strip ions of all their electrons. The number of atomic states retained in the partition function is a key quantity in determining if the EOS is complete at the conditions of interest. This assessment of completeness will of course strongly depend on the plasma conditions.

The final important aspect of opacity models that we wish to briefly discuss concerns line broadening. Photoabsorption lines in a plasma become broadened due to a number of factors, such as natural and Doppler broadening. Collisional broadening (often known as Stark broadening) is often a dominant broadening mechanism. This is broadening due to the plasma microfields caused by the motion of electrons and ions in a plasma. As before, line broadening is an important part of any opacity model, but any treatment of this effect has to be tractable for all ions and all conditions encountered when building an opacity table. While quite accurate line broadening models exist for one- and two-electron systems (Gomez *et al.*, 2022; Lee, 1988; Stambulchik *et al.*, 2019), the treatment of broadening for multi-electron systems often requires a number of approximations (Dimitrijevic and Konjevic, 1986, 1987).

C. Production of tables

Some of the earliest determinations of opacity for stellar physics started by the pioneering work of Seaton and

collaborators in the UK (Berrington *et al.*, 1987; Seaton, 1987). This development led to the “Opacity Project” (OP), which in time produced a set of opacity tables for stellar modeling that are still in common use today (Badnell *et al.*, 2005) At around the same time, OPAL (OPacity Astrophysical Library) opacity tables were published by Lawrence Livermore National Laboratory (LLNL) (Iglesias and Rogers, 1996). Both of these sets of tables were widely adopted by the stellar modeling community. Some early theoretical opacity work was performed at Los Alamos National Laboratory (Huebner and Barfield, 2014), starting from pioneering work by Mayer and later by Arthur Cox in the 1960s. In the 1990s, the LANL work resulted in “OPLIB” (OPacity LI-Brary) tables produced from the LEDCOP code (Magee *et al.*, 1995); in the mid 2010s the modern ATOMIC code was used to produce a new generation of opacity tables that improved the computation of opacities in a number of ways (Colgan *et al.*, 2016). All of these tables are available electronically in various formats. Other important opacity tables have been produced in France (the OPAS and SCO-RCG opacity efforts) (Blancard *et al.*, 2011; Mondet *et al.*, 2015; Pain, 2021), Israel (using the STAR code) (Krief *et al.*, 2016b, 2018a), and elsewhere. The comparison of opacity models with each other and with experiment (when available; see following sections) have been explored in dedicated workshops and conferences, for example Mendoza *et al.* (2018).

Many of these tables (primarily OP and OPAL) were quickly adopted by the solar modeling community and were useful in solving the Cepheid variability puzzle (Simon, 1982), where updated opacities helped resolve a previous disagreement between observation and modeling of Cepheid variables. Detailed comparisons between these two sets of tables have been made on numerous occasions, both on comparisons of individual opacities (Badnell *et al.*, 2005) and on how solar models that use these tables compare (Delahaye and Pinsonneault, 2006). Overall, the agreement between the various opacity tables was deemed to be “satisfactory” (within 5 %) for the accepted solar models of the late 1990s (Magee *et al.*, 1995). This paradigm was shaken when new determinations of the elemental solar abundances were reported (Asplund *et al.*, 2005), which implied that solar models disagreed with helioseismology observations. It was suggested (Serenelli *et al.*, 2009) that changes in the opacity of a few species (mainly Fe, Ne, O) could reconcile the solar models with the observations. Since then, intensive work has been performed in exploring some of the approximations made in the older opacity tables. The OP effort has focused on new and larger R-matrix calculations for important ions of Fe (Delahaye *et al.*, 2021; Nahar and Pradhan, 2016), primarily Fe^{17+} . At Los Alamos, the modern ATOMIC code was used to create a new generation of opacity tables for H through Zn (Colgan *et al.*, 2016). Although significant improvements in

many aspects of the calculations were incorporated, the resulting opacity tables did not significantly improve the comparison of solar modeling with helioseismology. The study of Krief *et al.* (2016a) showed that one possible way to produce larger opacities under solar conditions would be if line broadening effects were significantly increased – by several orders of magnitude. Opacity increases of $\approx 10\%$ near the solar convective zone and $\approx 2\%$ in the solar core have been attributed to ionic correlations (Krief *et al.*, 2018b). There is need for further investigation of such topics.

D. Experimental testing of calculated iron opacity

Errors in calculated solar opacity can arise from inaccuracies in both the abundance and calculated element opacities in a complex way. Therefore, it is important to investigate the accuracy of abundance and elemental opacities separately. The accuracy of elemental opacities can be best evaluated by examining the frequency-resolved (monochromatic) opacity. Since each spectral feature of the element opacity often relies on different physics and approximations, disagreements between model predictions and experimental data on a frequency-resolved basis can help identify which approximations are invalid and guide refinements in opacity theory. Here, we provide an overview of the fundamentals as well as previous investigations into measurements of solar-interior opacity. A more general and tutorial discussion of the experimental methods and a historical overview can be found elsewhere (Bailey *et al.*, 2009).

The material opacity κ_ν is related to its transmission T_ν by the following equation:

$$T_\nu = \frac{I_\nu - \epsilon_\nu}{B_\nu - \epsilon_\nu} = e^{-\kappa_\nu \rho L} \quad (33)$$

where B_ν is the backlight spectrum, I_ν is the backlight transmitted through the sample, ϵ_ν is a sum of the sample plasma emission and other backgrounds, and ρL is the areal density of the opacity sample. Thus, sample opacity κ_ν can be experimentally determined by heating the sample and accurately measuring the following quantities: the backlight with and without the heated sample (B_ν, I_ν), the plasma self-emission and background (ϵ_ν), and the sample areal density (ρL).

Reliable and useful opacity measurements require meeting several challenging criteria [Perry 1996, Bailey 2009]. First, the elements, conditions, and spectral ranges of interest must be identified. Second, a macroscopic opacity sample must be uniformly heated to the desired conditions to achieve local thermodynamic equilibrium. Third, the heated sample must be backlit with a bright and spectrally smooth radiation, B_ν , to accurately determine frequency-resolved absorption. It is critical that the backlight is significantly brighter than the

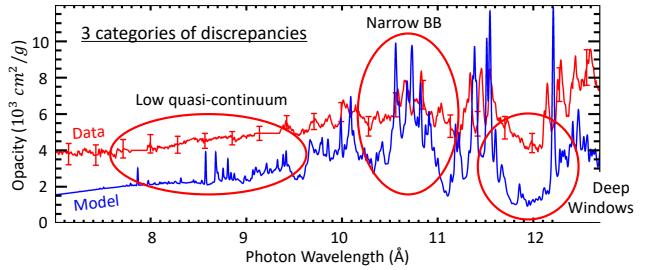


FIG. 12 Comparison of measured opacity (red) and calculated opacity (blue, Badnell and Seaton, 2003), at a temperature of 182 eV and an electron density of $3.1 \times 10^{22} \text{ e cm}^{-3}$. Three types of measured-vs-modeled opacity disagreements are shown: (i) a lower quasi-continuum, (ii) narrower bound-bound line features, and (iii) deeper opacity windows. These trends were observed in all opacity models compared in Bailey *et al.* (2015)

sample plasma emission or other background signals, ϵ_ν . Fourth, the absorption spectrum must be recorded with spectrometers that provide sufficient signals and spectral resolving power for the opacity study. Fifth, the condition of the heated sample ($T_e, n_e, \rho L$) must be diagnosed independently of the opacity in question. Simulated conditions are not appropriate for this purpose since their accuracy depends heavily on the accuracy of the opacity in question. Finally, the opacity spectrum and its uncertainty must be accurately determined from the measured $I_\nu, B_\nu, \epsilon_\nu$, and ρ_ν and their uncertainties.

After the revision of solar abundance in 2005 (Asplund *et al.*, 2005), solar models and helioseismology disagreed, and the accuracy of solar-interior opacity was called into question. This disagreement was primarily due to a significant reduction in solar opacity resulting from the reduced metallicity, and it was the greatest at the base of the solar convection zone (hereafter convection-zone base or CZB). Since then, solar abundance has been continuously revised, with the latest abundances determined by two groups being Z/X=0.0225 (Asplund *et al.*, 2021) and Z/X=0.0187 (Magg *et al.*, 2022), which still disagree with each other, leaving significant uncertainty in the solar abundance. For determining the accuracy of calculated solar opacity, the accuracy of calculated element opacities must also be experimentally scrutinized. Since Fe and O are the two dominant sources of opacity at the CZB, their opacities must be experimentally tested at CZB conditions (i.e., $T_e = 182 \text{ eV}, n_e = 9 \times 10^{22} \text{ e cm}^{-3}$). Unfortunately, previous experimental approaches were not suitable for this purpose because their backlights were not bright enough to mitigate the bright self-emission produced at these conditions.

Over the past two decades, experimental methods have been refined using one of the brightest x-ray sources, Z-pinch. In 2007, iron opacities were successfully measured at 150 eV and $7 \times 10^{21} \text{ e cm}^{-3}$, providing support for

the accuracy of calculated Fe opacities under these specific conditions. In 2015, iron opacities were measured at larger temperatures and densities. This study revealed a significant discrepancy between the calculated and measured opacities as the temperature and density approached the CZB conditions, as shown in Fig. 12. This is a significant concern since the experimental density was still approximately three times lower than the actual density at the CZB. Opacity models employ approximations to account for density effects such as line broadening and pressure ionization. If these approximations are incorrect, the disagreement between calculated and measured opacities could be even more pronounced at the densities found at the CZB.

The disagreement observed in the iron opacity was complex, as depicted in Fig. 12. The modeled opacities were lower in the quasi-continuum region at short wavelengths ($< 10 \text{ \AA}$), narrower in bound-bound lines, and deeper in the opacity valleys known as windows. These discrepancies are likely caused by various factors, but it is difficult to disentangle different sources of discrepancies based solely on the disagreement in iron opacity alone.

In 2019, a comprehensive study was conducted to measure opacities for Cr, Fe, and Ni at solar interior temperatures (Nagayama *et al.*, 2019). This systematic study helped narrow down hypotheses on the sources of disagreement by observing how the discrepancies change as a function of atomic number. The measurements revealed that measured opacities for all three elements had narrower line features than the model predictions, suggesting potentially inaccurate density effects in opacity models. Disagreement in the opacity windows was observed for Cr and Fe, while Ni did not show significant disagreement. At the experimental conditions used in the study, Cr and Fe were in an open L-shell configuration, whereas Ni was closer to a closed L-shell configuration. It is known that the calculations of both population and equation of state (EOS) are more challenging in open L-shell configurations. Therefore, the observed trend in window disagreement may suggest a connection to the difficulties associated with EOS calculations in such configurations.

One intriguing finding from this systematic study was the disagreement observed in the quasi-continuum region. Unlike the discrepancies observed in bound-bound linewidth and window opacities, no clear trend was observed in the quasi-continuum disagreement. The model and data agree well for Cr and Ni, but not for Fe, at high temperatures. Two possible explanations have been proposed for this unexpected behavior. The first hypothesis suggests the presence of missing physics in the opacity models that becomes significant under the conditions encountered by high-temperature Fe-opacity experiments. The second hypothesis proposes that the opacity experiments or analyses are flawed, but only for Fe at high temperatures. Until opacity experiments and theories are reconciled, neither possibility can be definitively ruled

out.

Despite numerous investigations, the discrepancies between experimental and theoretical iron opacities remain unresolved (Bailey *et al.*, 2015; Blancard *et al.*, 2016; Colgan *et al.*, 2016; Kilcrease *et al.*, 2015; Mancini, 2016; Nagayama *et al.*, 2016a, 2014a,b, 2016b, 2019, 2017; Nahar and Pradhan, 2016). On the experimental side, measurements of sample spatial gradients have been conducted and found to be negligible (Nagayama *et al.*, 2014a). Numerical tests have also shown that temporal gradients, self-emission, and the presence of tamping material do not have significant impact on the results (Nagayama *et al.*, 2014b, 2016b, 2017). Furthermore, the uncertainties in temperature and density diagnostics resulting from the choice of spectral model have been determined, and it has been found that these uncertainties are too small to account for the reported discrepancies (Nagayama *et al.*, 2016a). Additionally, background measurements have also been performed and found to be negligible (Dunham *et al.*, 2021).

The persistent differences between experiment (Bailey *et al.*, 2015; Nagayama *et al.*, 2019) and opacity models has led to a sustained effort to reconcile these discrepancies. The experimental efforts have been previously discussed. On the theoretical side, tests of some of the approximations used in the opacity tables have been conducted (Colgan *et al.*, 2016; Nahar and Pradhan, 2016; Blancard *et al.*, 2016; Delahaye *et al.*, 2021), particularly in assessing the convergence of opacity models with respect to the number of configurations included in the calculations. However, no significant change in the final opacity has been reported from such studies.

Several groups have postulated new physical effects that could produce more opacity contributions than in previous works. Of particular note is the possibility of two-photon opacity contributions, where an atom or ion could simultaneously absorb two photons (possibly of different frequency) (More *et al.*, 2020, 2017). All current opacity tables omit such contributions, as they are believed to be a minor contribution to the overall opacity. Preliminary studies by More and co-workers (More *et al.*, 2020, 2017) have suggested this could be a factor in the comparison of theory with the Z measurements, but other studies (Iglesias, 2015; Kruse and Iglesias, 2019, 2021; Pain, 2018) find that such a contribution should be small.

Another intriguing study has reported that transient spatial localization may result in increased opacities (Liu *et al.*, 2018; Zeng *et al.*, 2022). This is a density-dependent effect where the plasma effects localize the continuum wavefunctions near the absorbing ion, potentially resulting in increased broadening and bound-free cross sections. While the work of (Zeng *et al.*, 2022) suggests that this effect is significant, other studies cast doubt on this claim (Iglesias, 2023).

Opacity theorists and experimentalists continue to col-

laborate closely to address discrepancies and uncertainties in the calculated solar RMO. They carefully examine each other's work, seeking to resolve the discrepancies and accurately quantify the uncertainty associated with calculated solar opacity.

E. Future opacity work

At the time of writing, the continuing disagreement between the Z Fe experiments ([Bailey et al., 2015](#); [Nagayama et al., 2019](#)) and theoretical models is by far the biggest concern of the opacity community. New theoretical efforts, while intriguing, have not yet resolved these differences. New experimental results, particularly the opacity-on-NIF campaign ([Opachich et al., 2022](#); [Perry et al., 2020](#)), are eagerly anticipated to shed some light on the current impasse. New approaches are also being developed for opacity measurements using the APOLLON laser at LULI (France) and on the GEKKO XII laser (Japan) ([Fujioka et al., 2005](#)). There has been a suggestion for measuring heat conductivity under thermodynamic conditions resembling those found deeper in the radiative zone, using hydro-carbon foams doped to imitate the solar composition ([Krief et al., 2018b](#)). Such progress will have implications for solar modeling, as the combination of opacity and elemental abundances is key to resolving the significant discrepancy between helioseismology and solar models. As previously noted, the abundance of solar elements is also the subject of renewed scrutiny ([Asplund et al., 2021](#); [Magg et al., 2022](#)).

XIV. Experimental facilities for solar fusion studies

Improved experimental facilities have led to significant progress in nuclear astrophysics research in recent years, and future instrumental developments will likely continue to push boundaries in the next decade. Key goals driving technological developments include, among others, tests of the weak interactions and of solar properties that make use of high precision solar neutrino measurements, helioseismological mappings of the core metallicity, and detailed solar modeling. There is also a host of related open questions in different areas, such as Big Bang nucleosynthesis, red giant evolution, the evolution of supernova progenitors, and a variety of transient explosive phenomena in astrophysics where a quantitative understanding of the nuclear physics is essential. Due to the small cross sections of low-energy charged-particle reactions, experiments must be designed to achieve signal rates significantly lower than the background rates from cosmic rays, natural radioactivity in the laboratory, and induced activity from beam interactions with target impurities. In the following, we outline current experiences and future facilities—either in development or planned—that will

pave the way for possible major breakthroughs in the coming years.

A. Above-ground facilities

Above-ground facilities always play a crucial role in investigating nuclear reactions of astrophysical interest. Various methods are applied to minimize background. A prevalent approach involves employing passive shielding around the detection area, typically utilizing a layered configuration of lead, copper, and polyethylene. This setup helps to decrease unwanted signals and neutron background in detectors with relatively small capacities. However, there are additional techniques that can be utilized to further diminish background interference, enabling measurements at energy levels closer to those pertinent to astrophysics.

1. LENA

The Laboratory for Experimental Nuclear Astrophysics (LENA) is part of the Triangle Universities Nuclear Laboratory, located on the campus of Duke University, in North Carolina, US. LENA features two accelerators: a 230-keV ECR accelerator that produces the world's most intense low-energy proton beams ([Cooper et al., 2018](#)) and a 2-MV Singletron from High Voltage Engineering Europa B.V. ([Shornikov et al., 2023](#)). These machines deliver ion beams with remarkable intensity—up to 20 mA DC proton current at lower energies and 2 mA at several MeV—along with advanced beam pulsing capabilities. Each accelerator has dedicated transport systems and control mechanisms for simultaneous operation. A key component of LENA is the γ -ray coincidence spectrometer, which allows measurements with sensitivities comparable to underground facilities ([Buckner et al., 2015](#)). The detector setup includes a 130% p-type HPGe detector, HPGe clover detectors, and a 16-segment NaI(Tl) annulus, along with the APEX detector featuring 24 position-sensitive NaI(Tl) bars for enhanced measurements. Recently, LENA added a cosmic-ray veto system with nine plastic scintillators to reduce cosmic-ray muon background, conveniently positioned due to its wheeled assembly.

2. NSL

The Nuclear Science Laboratory (NSL) at the University of Notre Dame, hosts a number of low energy accelerators and the TriSOL radioactive beam facility to maintain a rigorous experimental program in nuclear astrophysics, nuclear structure physics, and fundamental symmetries as well as in a broad range of nuclear physics and nuclear chemistry applications. The 5MV single

ended pelletron provides high intensity proton, helium, and $A \leq 40$ heavy ion beams for nuclear astrophysics related experiments in forward kinematics using solid and gas target technologies and in inverse kinematic using the St. George recoil separator. The 10 MV FN pelletron tandem is the main machine for nuclear structure physics experiments, but also serves that AMS program for analysing long-lived radioactivities in terrestrial and meteoritic samples. The FN is also used for indirect measurements of reactions of astrophysical interest and serves as a driver for the production of light radioactive isotopes beams ($A \leq 40$) for nuclear astrophysics and fundamental symmetry studies.

B. Underground Facilities

Underground facilities are essential to push the boundary of direct measurements towards the lowest energies of astrophysical interest (10s-100s keV). In particular, capture-reaction measurements leading to neutron- or γ -ray emissions are severely hampered in surface laboratories because of the overwhelming background associated with cosmic rays. This background can be suppressed by many orders of magnitude by exploiting the natural shielding provided by the rock overburden in underground sites. The improvements possible with this strategy have been demonstrated by pioneering work at the Laboratory for Underground Nuclear Astrophysics (LUNA) at INFN-LNGS, first with a 50-kV accelerator (LUNA I) and then with a 400-kV one (LUNA II) still in operation today (Aliotta *et al.*, 2022a). Since the last Solar Fusion review, a new accelerator has been installed at INFN-LNGS, and other underground laboratories have become operational in China (JUNA), in the US (CASPAR), and in Germany (Felsenkeller). These are briefly presented below.

1. Bellotti Ion Beam Facility

The INFN-LNGS has recently expanded its accelerator capabilities with the installation of a new 3.5 MV SingletronTM machine designed and set up by High Voltage Engineering Europe (HVEE) (Sen *et al.*, 2019), see Fig. 13. The 3.5 MV machine is equipped with two independent beam lines which can be operated with solid and gas target systems. Acceptance tests at HVEE demonstrated that the machine can deliver intense proton, helium and carbon beams (1, 0.5 and 0.15 mA, respectively) with well defined energy resolution (0.01% of TV) and stability (0.001% h^{-1} of TV) (Di Leva, 2020 - private comm.). The 1.4km of rock overburden (ca. 3800 m.w.e.) at LNGS affords significant background reductions for both γ -ray, neutron-, and charged-particles detection (see (Aliotta *et al.*, 2022a) for a recent review). A first exper-



FIG. 13 Photo of the 3.5 MV SingletronTM accelerator recently installed at LNGS (credit: Matthias Junker).

imental proposal presented by the LUNA-Collaboration focuses on measurements of the reactions $^{14}\text{N}(\text{p},\gamma)^{15}\text{O}$, $^{12}\text{C}+^{12}\text{C}$, $^{13}\text{C}(\alpha,\text{n})^{16}\text{O}$ and $^{22}\text{Ne}(\alpha,\text{n})^{25}\text{Mg}$, the latter being in the context of the ERC Starting Grant *SHADES*. The 3.5MV accelerator is now part of the Bellotti Ion Beam Facility (Junker *et al.*, 2023) and open to external users.¹³

2. JUNA

The Jinping laboratory for Underground Nuclear Astrophysics (JUNA) is part of the China Jinping Underground Laboratory (CJPL), established on the site of hydropower plants in the Jinping mountain, Sichuan, China (Liu *et al.*, 2022b). The laboratory, located near the middle of a traffic tunnel, is shielded by 2400 m of mainly marble overburden (6720 m.w.e.). In December 2020, the JUNA collaboration installed a 400 kV accelerator capable of delivering H^+ and He^+ beams with intensities of up to 10 particle mA, and a He^{2+} beam with an intensity of up to 1 particle mA. Some important results have already been published. For example, the lowest energy study to date of the $^{19}\text{F}(\text{p},\gamma)^{20}\text{Na}$ reaction advocates a breakout from the CNO cycle as a possible route for calcium production in population III stars (Zhang *et al.*, 2022).

3. CASPAR

The Compact Accelerator System for Performing Astrophysical Research (CASPAR) laboratory is the only US-based deep underground accelerator and is operated by a collaboration of the University of Notre Dame and the South Dakota School of Mines and Technology (Robertson *et al.*, 2016). The accelerator system has been

¹³ <https://www.lngs.infn.it/en/pagine/bellotti-facility-en>

fully operational since 2018 and is located 1480 m below surface at the Sanford Underground Research Facility (SURF)¹⁴ in Lead, South Dakota, formerly the Homestake gold mine. The rock overburden results in a 4300 m.w.e shielding effect, significantly decreasing cosmic-ray induced background with a muon flux level of 4×10^{-9} cm⁻² s⁻¹. The residual neutron flux consists of primarily low-energy (<10 MeV) neutrons generated by (α, n) reactions induced by the decay of naturally occurring uranium and thorium radionuclides in the surrounding rock, and is generally of the order of 10⁻⁶ neutrons cm⁻² s⁻¹ (Mei and Hime, 2006; Mei *et al.*, 2009). The CASPAR accelerator is a 1 MV Van de Graff style JN accelerator with a 150 to 1100 kV operational range, well suited for overlap with higher energy measurements. The accelerator provides proton and α beams with up to ~ 250 μ A on target. The scientific program continues to explore stellar neutron sources and expands the present studies into the magnesium range probing α -capture reactions on ²⁴Mg, ²⁵Mg, and ²⁶Mg isotopes. A new program has been initiated to explore the endpoint of nova nucleosynthesis, studying proton capture reactions in the Ar to Fe range. CASPAR is well suited for these measurements, but will be complemented by a new low energy machine, presently under development at Notre Dame.

4. Felsenkeller

The Felsenkeller laboratory is hosted within a network of tunnels, excavated into hornblende monzonite rock, previously used as a cool storage place for the homonymous nearby brewery. Jointly funded and built by TU Dresden and HZDR, the laboratory operates a 5 MV Pelletron-type accelerator built by NEC. With the addition of a RF ion source in the high-voltage terminal, the accelerator can be operated in tandem mode or as a single-ended machine and is capable of delivering proton, α -particle, and carbon beams with currents up to several tens of μ A. Unlike the other sites mentioned in this section, Felsenkeller is a shallow underground laboratory, with a rock overburden of only 140 m.w.e. Such a depth is sufficient to shield all components of cosmic-ray induced radiation, except muons, for which further active shielding is generally required (Ludwig *et al.*, 2019). Detailed studies of muon, neutron, and γ backgrounds have been carried out (Grieger *et al.*, 2020; Ludwig *et al.*, 2019; Szucs *et al.*, 2019a; Turkat *et al.*, 2023), and the ¹²C(p, γ)¹³N reaction has been studied (Skowronski *et al.*, 2023b). The further scientific program of Felsenkeller foresees the study of several reactions of astrophysical interest, including ³He(α, γ)⁷Be and ¹²C(α, γ)¹⁶O.

C. Indirect methods

Indirect methods are also pivotal in nuclear astrophysics due to the challenges associated with directly measuring reactions at extremely low energies. They can have systematic uncertainties that are different from those of direct measurements, and they provide supplementary information that can constrain R-matrix and other models used in the extrapolation of data from direct measurements. For the present review, it is worth further exploring here the ANC (asymptotic normalization coefficient) and THM (Trojan Horse) methods. For a detailed description of the methods, refer to (Tribble *et al.*, (2014; Tumino *et al.*, 2021).

The ANC method determines the zero-energy cross-section for radiative-capture reactions by exploiting their peripheral nature. This means the reaction depends mostly on the long-distance behavior of the wave function. The ANC is extracted from transfer reactions using the DWBA, with uncertainties primarily arising from the optical model description, still much smaller than those affecting spectroscopic factors. While the method is focused to zero energy, it complements direct measurements and can be applied to loosely bound nuclei. The method has been applied to the ³He(α, γ)⁷Be (section VI.C).

The THM is a powerful technique for indirectly measuring astrophysical S(E) factors of reactions involving charged particles. By studying a related reaction with a spectator particle, the THM allows for the extraction of the desired cross section without the need for extrapolation. The THM relies on the assumption of quasifree kinematics, where the spectator particle has a minimal effect on the reaction of interest. The method requires careful selection of beam energy and momentum transfer to ensure that the Coulomb barrier is overcome and the spectator particle remains relatively undisturbed. Since Solar Fusion II, significant advancements have been made, focused on improving the model's accuracy and assessing its systematic uncertainties (Tribble *et al.*, (2014; Tumino *et al.*, 2021). For reactions dominated by broad resonances, the modified R-matrix approach (La Cognata *et al.*, 2015b; Trippella and Cognata, 2017) allows for the incorporation of half-off-energy-shell and energy resolution effects within a well-established framework. This method enables multi-channel descriptions of reactions, such as in the ¹²C+¹²C fusion studies (Tumino *et al.*, 2018), and includes a DWBA-based normalization procedure that does not rely on direct data (La Cognata *et al.*, (2009). The modified R-matrix framework has been applied to the ¹⁵N(p, α)¹²C (section XI.A.3), ¹⁹F(p, α)¹⁶O (section XI.C.4), and ²³Na(p, α)²⁰Ne (section XI.D.5) reactions. For narrow resonance reactions, a simplified approach has been introduced (La Cognata *et al.*, 2022) to deduce resonance strengths. This method reduces systematic errors from normalization and the-

¹⁴ See <http://www.sanfordlab.org>

ory to the percent level through multi-resonance normalization and covariance in error propagation. The narrow resonance approach has been used in analyzing the $^{17}\text{O}(\text{p},\gamma)^{18}\text{F}$ (section XI.B.3), $^{17}\text{O}(\text{p},\alpha)^{14}\text{N}$ (section XI.B.4), and $^{18}\text{O}(\text{p},\alpha)^{15}\text{N}$ (section XI.C.2) reactions.

D. Plasma facilities

1. NIF/OMEGA

Several efforts have recently been devoted to studying nuclear reactions in a plasma environment reminiscent of stellar conditions (Gatu Johnson *et al.*, 2017) using the two large laser facilities: OMEGA at the Laboratory for Laser Energetics, University of Rochester (Boehly *et al.*, 1997), and the National Ignition Facility (NIF) at Lawrence Livermore National Laboratory (Wontergem et al., 2016). Both facilities use high-power lasers to implode spherical capsules containing reactants of interest to high temperatures (of order 1-20 keV) and densities (up to 10^3 g cm^{-3} in the extreme case of implosions with a fuel ice layer (Abu-Shawareb *et al.*, 2022)). The OMEGA laser has the capability of delivering 30 kJ of laser energy divided between 60 laser beams to the target; the NIF can deliver up to 2.0 MJ using 192 laser beams, and can thus be used to implode larger amounts of materials to more extreme conditions compared to OMEGA. Nuclear experiments at these two facilities are enabled by an extensive suite of nuclear diagnostics, originally developed to do inertial confinement fusion experiments (Cerjan *et al.*, 2018). Initial results have been obtained on the $^2\text{H}(\text{p},\gamma)^3\text{He}$ reaction (Mohamed *et al.*, 2022; Zylstra *et al.*, 2020), the $^3\text{He}(^3\text{H}, \gamma)^6\text{Li}$ reaction (Zylstra *et al.*, 2016), and the $^3\text{H}(^3\text{H},2\text{n})\alpha$ reaction (Casey *et al.*, 2017). Proton spectra have also been measured for the $^3\text{He}(^3\text{He},2\text{p})\alpha$ reaction (Zylstra *et al.*, 2017). In addition to low-Z reaction nuclear S-factor studies, the NIF and OMEGA facilities are also promising for studies of plasma effects on nuclear reactions, including screening (Aliotta and Langanke, 2022; Casey *et al.*, 2023); for studies of charged-particle-induced reactions (Wiescher *et al.*, 2022); and, thanks to the high neutron fluxes achievable (up to $5 \times 10^{27} \text{ neutrons cm}^{-2} \text{ s}^{-1}$ (Abu-Shawareb *et al.*, 2022)), for studying reactions on excited states (Thompson, 2022). However, there are challenges remaining to be addressed in fully developing this new platform for nuclear experiments, including the impact on the results of rapid gradients in space and time (Crilly *et al.*, 2022).

2. PANDORA

PANDORA (Fig. 14) is a device conceived for multidisciplinary studies, including many of astrophysical interest (Mascali *et al.*, 2022). Two of its main objectives are: (a)

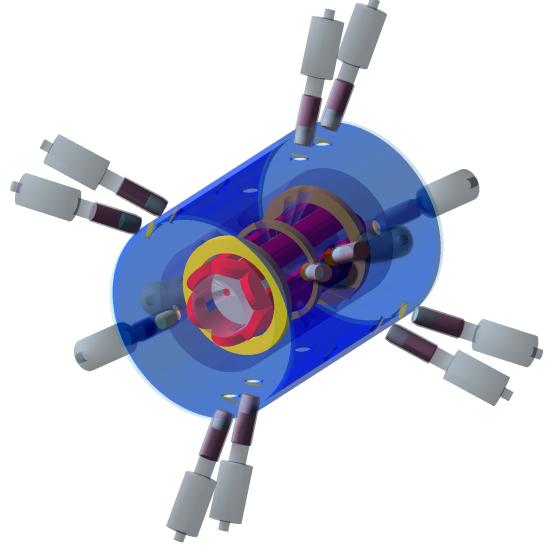


FIG. 14 Schematic drawing of the PANDORA trap with holes to house the HPGe detectors. (credit: Domenico Santonocito).

to perform the first measurements of β decays in plasmas of astrophysical relevance, in order to verify the results obtained in storage rings with ^{187}Re (a lifetime reduction by 9 orders of magnitude), and (b) to measure the opacities of plasmas of astrophysical interest (kilonova ejecta). PANDORA will mainly consist of three subsystems:

- an innovative superconducting magnetic plasma trap, able to produce and confine plasmas with electron-ion density up to 10^{13} cm^{-3} and electron temperature of $T_e \sim 0.1 - 30 \text{ keV}$;
- an advanced plasma multi-diagnostic system, consisting of a set of non-invasive diagnostic tools capable of operating simultaneously, for non-intrusive monitoring of the thermodynamic plasma properties and parameters;
- an array of 14 HPGe (High-purity Germanium) detectors for γ -ray spectroscopy, surrounding the plasma trap.

The next development steps will involve: preliminary numerical simulation studies to assess the possibility of measuring opacities in astrophysical plasmas (relevant for kilonova ejecta); study of expected abundances and constraints in AGB stars for some (key) nuclides of interest; experimental investigation of magnetic confinement and turbulence in plasma, using an existing trap at ATOMKI-Debrecen; definition of the analysis algorithm for X-ray imaging and spatially resolved spectroscopy.

Initial physics cases to be investigated include ^{176}Lu (a potential cosmo-chronometer), ^{134}Cs (to reproduce adequately the observed abundance ratio of the two s-only

isotopes, ^{134}Ba and ^{136}Ba) and ^{94}Nb (to solve the puzzle about the exact contribution of s-processing to ^{94}Mo).

3. Sandia Z

The Sandia National Laboratories (SNL) Z machine is the world's largest pulsed power accelerator ([Sinars et al., 2020](#)). It charges 22 MJ of electrical energy into a capacitor and then discharges it all at once (over approximately 100 ns) into targets ranging from millimeters to centimeters in size. This allows it to reach a peak electrical power of 80 terawatts, which means that, for that instance, it generates 15 times the electrical power generated by all of the world's power plants combined. By concentrating this massive power into small targets, the Z machine can convert them into a state of high energy density (HED), which refers to an extreme state of matter with a pressure exceeding 10^6 times atmospheric pressure.

The Z machine has been used for a wide range of HED science, including inertial confinement fusion, dynamic material properties, and laboratory astrophysics. In particular, the Sandia Z machine can convert this electrical power into X-ray power using a scheme called the Z-pinch dynamic hohlraum ([Rochau et al., 2014](#)). This is the most energetic X-ray source on earth and has been used by academic collaborations such as the Wooton Center for Astrophysical Plasma Properties (WCAPP) and the Center for Laboratory Astrophysics. For the stellar opacity project (see Sect. [XIII.D](#)), they heat iron or oxygen to the conditions relevant to the base of the convection zone and measure their frequency-resolved opacities to test hypotheses for the solar opacity-abundance problem ([Bailey et al., 2015, 2009; Basu and Antia, 2008; Nagayama et al., 2019](#)). For the accretion disk projects, they heat silicon, iron, and neon photoionized plasmas to conditions similar to those at black hole accretion disks to experimentally test the validity of calculated atomic data, kinetics, and spectral formation ([Loisel et al., 2017](#)). This is important for accurately interpreting high-resolution data from upcoming X-ray space telescopes such as XRISM. For the white dwarf photosphere project, they reproduce its photosphere conditions and validate spectroscopic methods used for understanding the age of the universe as well as constraining its evolutionary paths and supernova remnants ([Falcon et al., 2015; Montgomery et al., 2015; Schaeuble et al., 2021, 2019](#)). These experiments reach temperatures ranging from 10^4 to 2×10^6 K and densities ranging from 10^{16} to 10^{23} electrons cm^{-3} and are performed simultaneously by placing their samples at different locations from the Z-pinch dynamic hohlraum.

E. Storage rings for nuclear astrophysics studies

1. CRYRING

A main limitation of experimental studies involving unstable nuclei arises from the difficulty of producing radioactive ion beams (RIBs) of adequate intensity and purity at the energies of interest for astrophysical applications. Storage rings ([Steck and Litvinov, 2020](#)) offer a key advantage over traditional ISOL and in-flight approaches as they allow for storing and re-circulating radioactive ions over and over again, thus allowing for multiple interactions of un-reacted beam particles with an in-ring target. Beam re-circulation significantly improves the quality of the beam as it results in orders-of-magnitude increased intensity (typical boosting factors of $\sim 10^5$), limited only by the duty cycle of the measurement ([Bruno et al., 2023](#)), and in improved purity of the stored beam, as only beam particles with the right mass-to-charge ratio will survive in the beam orbit.

One of the main technical challenges of storage rings comes from the requirement of ultra high vacuum (UHV) conditions ($< 10^{-10}$ mbar) necessary to guarantee that ions can be recirculated with minimal losses. This, in turn, translates into a requirement for sufficiently thin targets (i.e. typically $\leq 10^{11} - 10^{14}$ atoms cm^{-2}) to minimise beam ion losses through scattering or electron capture (see ([Bruno et al., 2023](#)) for further details).

A dedicated low-energy storage ring, CRYRING ([Lestinsky et al., 2016](#)), inherited from Stockholm University, was installed at GSI in 2016, as a Swedish in-kind contribution to the FAIR facility. The facility allows nuclear astrophysics studies to be done with radioactive nuclei. After its initial recommissioning ([Herfurth et al., 2018](#)), CRYRING now serves as a low-energy extension for Experimental Storage Ring (ESR) beams, as well as a standalone machine with a local ion source ([Geithner et al., 2017](#)). After in-flight production in the Fragment Recoil Separator (FRS) at relativistic energies, rare ions can now be cooled, post-decelerated and stored in the full range down to about 100 keV/ u . A study is underway to reach lower energy by means of a transverse low-energy beamline from the FISIC project ([Glorius and Bruno, 2023; Schury et al., 2020](#)). CRYRING is now equipped with an in-ring micro-droplet gas target, whose low density ($\leq 10^{14}$ atoms cm^{-2}) ensures minimal beam energy loss and straggling through the target, a unique advantage for charged particle spectroscopy at storage rings compared to standard techniques.

In addition, a new detection chamber (Fig. [15](#)), CARME (CRYRING Array for Reaction MEasurements), was recently installed at CRYRING and fully commissioned in early 2022. Specifically designed to fulfil requirements of stored beam experiments and operating at a pressure of 10^{-12} mbar, CARME can host up to eight double-sided-silicon-strip detectors with excellent

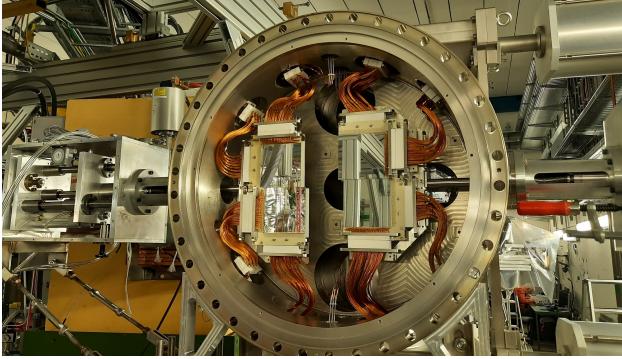


FIG. 15 Photo of the CARME chamber recently installed at CRYRING. Four silicon strip detectors are visible at the centre of the chamber (credits: Carlo Bruno).

angular and energy resolution.

Finally, the upcoming installation of the FISIC transverse beam line (Schury *et al.*, 2020) will allow to intercept the beam stored in CRYRING, thus opening up unprecedented opportunities for crossed-beam experiments. An exciting application may be the study of nuclear reactions between ion beams, i.e. unaffected by the electron screening, directly at the energies of astrophysical interest. Initial proposals for the study of nuclear astrophysics reactions have already been approved.

XV. Closing remarks

A. Summary

This review summarizes the significant progress that has been made over the last decade in understanding the nuclear reactions that govern energy production in solar and stellar hydrogen burning. It also describes some of the plasma and atomic physics that effects the solar environment in which these reactions take place, as well as the diagnostics tools – solar neutrinos and helioseismology – by which we can probe that environment.

For S_{11} and S_{12} , cross sections have been derived from first principles, using both potential theory and effective field theory. In the case of S_{12} , new experimental data agree well with the theoretical predictions. S_{hep} , which governs a minor branch of solar fusion, must also be taken from theory.

The S-factors for other pp-chain reactions are based on laboratory measurements. In some cases, including notably S_{17} , most of the studies reviewed here are relatively far from the solar Gamow peak energies, underlining the need for theoretical and experimental tools to bridge the gap.

The nuclear physics of the pp chain remains a significant uncertainty in SSM predictions of the individual neutrino fluxes - pp, pep, hep, ${}^7\text{Be}$, ${}^8\text{B}$, and CNO. The nuclear errors are comparable to, and in the case of the

hep neutrinos greatly exceed, the “environmental” errors that are generated by other uncertainties in the SSM (Orebi Gann *et al.*, 2021).

For several reactions including S_{33} and S_{14} , the SF III recommended values are less precise than those of SF II. This fact should not be taken as a sign of regression, but instead as indicative of a more cautious treatment of uncertainties. In some of the error estimates made in this review, we have been mindful of past lessons on the potential impact of unidentified systematics.

The status of our understanding of radiative opacities has been reviewed here, the first time the subject has been included in the Solar Fusion series. The debate generated by the solar composition problem has brought renewed attention to the complicated interplay between opacities and composition in the SSM. Given new experimental opportunities, especially luminous X-ray sources coupled with plasma targets, there is the expectation of rapid progress in this field over the next decade.

There has been only very limited progress on electron screening since SF II. This field has important practical implications not only for the Sun but also for laboratory experiments, and may merit additional attention in the coming decade. Progress may depend on achieving a better understanding of low-energy stopping powers.

B. Recommended values for S-factors and their derivatives

The SF III recommendations for nuclear reaction S-factors and their derivatives are presented in Table I of the Introduction. In the sections that follow the Introduction, we presented the details leading to these results and identified opportunities for future work, which are summarized directly below.

C. General recommendations for future work

For all of the solar fusion reactions, obtaining high-precision experimental data in Gamow window remains an elusive goal, one that may not be reached in the foreseeable future. However, such direct measures are just one of many avenues for improving our understanding of the nuclear physics of the Sun. The recommendations below are in most cases summaries of more detailed discussions presented in this review, and we refer readers to the relevant sections for additional details.

1. For all of the experimentally accessible pp-chain reactions and for S_{14} which determines the CNO cycle rate, there is a need for at least one experimental data set spanning the entire energy range, encompassing the more limited data sets that currently exist. This will provide an important crosscheck on the normalizations of those data sets. Experimentalists can now tackle this challenge due to the new

- generation of underground accelerators that have the dynamic range needed for such studies.
2. We recommended continued effort on reaction theory relevant to pp-chain reactions. The needs go beyond rates to include both astrophysical (e.g., precise neutrino spectra) and laboratory (e.g., γ -ray angular distributions) observables.
 3. We urge the community to set an ambitious goal for the precision of future S-factor measurements: reduce the nuclear physics uncertainties to a factor of two below current SSM “environmental” uncertainties, so that they are no longer a significant contributor to SSM neutrino flux uncertainties. With the exception of the hep neutrinos, this goal could be reached by a future reduction of a factor \sim two, typically, in the individual S-factor uncertainties ([Orebi Gann et al., 2021](#)). As solar neutrino data are an important input into global neutrino oscillation analyses, nuclear uncertainties will continue to feed into that program until this goal is achieved. Improvements in S_{14} and other CNO cycle S-factors are needed, as the current 8.4% uncertainty would be a limiting factor in extracting the solar core’s metallicity from a future large-volume CN neutrino experiment. The progress reported here on S_{17} , with the uncertainty reduced significantly from [SF II](#) to 3.4%, should be continued, given the current and future ^8B neutrino programs of Super-K and Hyper-K: these neutrinos are our best solar core thermometer. Finally, ongoing Super-K efforts to detect hep neutrinos provide strong motivation for improving estimates of S_{hep} and the associated neutrino spectrum. (Uncertainties in the high-energy tail of the ^8B spectrum must also be reduced, as discussed in Section [II](#).)
 4. New experimental methods are becoming available that could help us better understand electron screening in terrestrial and solar reactions. Plasma conditions resembling those of the Sun can now be produced in the laboratory. Reactions can be studied in rings, using ions stripped of their atomic shells. The level of theory activity in this field has declined over the past decade, so we hope experimental progress will help renew interest.
 5. In order to break the degeneracy between solar abundances and solar opacities, efforts should be made to improve our understanding of radiative opacities for the Sun’s principal metals. A high priority is the resolution of the current discrepancy between measured and modeled iron opacities. Solar physicists need access to state-of-the-art, open-source opacity codes.
 6. We recommend that resources be available to extend the use of current and next-generation dark matter detectors to solar neutrino detection. This includes designing new detectors so that they are highly capable for both applications. Some of the high-priority goals of solar neutrino physics – such as a 1% measurement of the pp neutrinos to check the luminosity constraint – might be achieved with a dual-purpose detector.
 7. The Solar Fusion program to periodically review the nuclear physics of the Sun and other hydrogen-burning stars should be continued, with the separation between studies being no more than ten years. The format should continue to be open, giving all researchers working in the field an opportunity to contribute, and broadly international.
 8. Future Solar Fusion studies should continue to strengthen the connections between this community and others interested in main sequence stars. SF III has placed increased emphasis on hydrogen burning at higher temperatures, plasma physics, opacities, and asteroseismology. However, with observations now being made of solar-like stars in their formation stages, we anticipate additional connections developing, well beyond the limits of this study.

D. Outlook

Six decades after the initial observations of solar neutrinos by Ray Davis ([Davis, 2003](#)), significant advances have been made in our understanding of the SSM and indeed of hydrogen-burning stars in general. The driver of this progress has been experiment: the observations made of all of the principal solar neutrino sources; our improved understanding of the flavor physics of those neutrinos; the advances made in helioseismology, a second quantitative probe of the solar interior; and our improved knowledge of the input microphysics of the SSM, including nuclear cross sections and opacities. Despite this progress, improvements are still needed. In several cases the dominant uncertainty in our solar neutrino flux predictions stems from the limited precision of the nuclear S-factors. Similarly, uncertainties remain in our understanding of radiative transport in the Sun, including the continuing debate about the Sun’s primordial composition. As changes in composition often can be mimicked by changes in radiative opacities, our imperfect knowledge of the latter has slowed resolution of the solar composition problem.

The prospects for substantial progress in the next decade are bright. Whereas at the time of [SF II](#), there was only one underground ion accelerator worldwide, now there are five, opening up possibilities for independently

cross-checking important nuclear data. The low background that can be achieved underground has led to the first measurements of reactions in the Gamow window. Other new experimental capabilities have come on line that are advancing our understanding of solar atomic physics, including low-energy storage rings and new plasma facilities, the latter made possible in part by advances in high-power lasers. The field is approaching the point where opacity measurements can be made under conditions closely approximating those of the solar interior.

Theory has progressed as well, including advances in neutrino flavor physics that have made neutrinos once again a precise probe of solar physics. Techniques with a firmer footing in first principles, such as effective field theory and lattice QCD, are being employed in the extraction of pp chain S-factors. Conventional approaches like the R-matrix can now be applied to complex nuclear systems where the inclusion of multiple reaction channels is important.

On the observational side, several very large neutrino detectors are under construction, including Hyper-Kamiokande, JUNO, and DUNE. Certain direct-detection dark matter experiments have reached a sensitivity where nuclear recoils due to the coherent solar neutrino scattering have become a background, leading to discussions of how such detectors might be optimized for solar neutrino detection (Essig *et al.*, 2018). This underscores the remarkable radiopurity of these underground detectors. Finally, the helioseismic studies that proved so important in establishing the credibility of the SSM are now being extended to other stars – a major reason SF III was undertaken at this time. Asteroseismic studies have the potential to tell us what is special and what is common about the Sun as a star.

Given current excitement about the emerging field of multi-messenger astrophysics, it is perhaps fitting to remember that the Sun was the prototype for this field. An array of physical measurements of the Sun – its mass, radius, luminosity, abundances – were combined with neutrino flux and helioseismic measurements to constrain and then test the SSM. The result was the discovery of neutrino mass and flavor mixing.

This success was a multi-disciplinary effort that has not yet run its course. We still have uncertainties in our characterization of the Sun’s nuclear, atomic, and weak interaction physics that, if improved, would make the Sun an even more powerful laboratory. As our best known and nearest star, the Sun is the test of our understanding of the structure and evolution of other main-sequence stars. If the past is any indicator of the future, the efforts we make over the next decade to make the Sun an even more precisely calibrated reference will yield a rich return in our understanding of physics.

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A. Gallium Neutrino Source Cross Sections

Here we describe in more detail the constraint that the electron capture rate for ^{71}Ge places on the ^{51}Cr and ^{37}Ar neutrino source cross sections for ^{71}Ga , expanding on the discussion of Section II.A.2. We discuss both the early treatment of Bahcall and the more recent analysis of Elliott *et al.* (2024, 2023), performed in support of the BEST experiment

Following Bahcall (Bahcall, 1978), the electron capture rate can be written in terms of a dimensionless phase space factor f_{EC} :

$$f_{\text{EC}} = 2\pi^2 \left(\frac{\hbar}{mc} \right)^3 \left(\frac{q}{mc^2} \right)^2 |\psi(R)|^2, \quad (\text{A1})$$

where q is the energy available to the neutrino and we have omitted atomic shell effects for brevity. The last factor is the electron density at the nucleus within the radius R . Multiplying this by the experimentally determined half-life $t_{1/2}$ yields the traditional ft value needed to calculate the neutrino cross section. Bahcall (1978) defines a characteristic scale σ_0 for neutrino cross sections on ^{71}Ga in the form

$$\sigma_0 = 1.2429 \times 10^{-47} \left(\sum_i q_i^2 g_i^2 \right)^{-1} \quad (\text{A2})$$

$$= 8.611 \times 10^{-46} \text{ cm}^2 \quad (\text{A3})$$

where $g_i = \psi(R)$. Bahcall (1997) reports the number given above and states that “it is about 0.5% less than the value 8.8012×10^{-46} I have used since 1984.” Since these numbers differ by 2.2%, there is clearly an error in one of the 3 numbers.

Recently Elliott *et al.* (2024, 2023) have recast the derivations with fundamental constants replacing numerical values and correction terms explicitly stated. They conclude that the ground-state cross section is 2.5% lower, and the excited-state contributions somewhat larger, than values used in the anomaly analysis. The adjusted net anomaly is minimally reduced.

They find the following expression for the neutrino cross section on ^{71}Ga leading to the ground state of ^{71}Ge :

$$\sigma_{\text{gs}} = \frac{G_F^2 \cos^2 \theta_C g_A^2}{\pi} p_e E_e \mathcal{F}(Z_f, E_e) B_{\text{GT}}^{(\nu,e)}(\text{gs}) [1 + g_{v,b}(\nu,e)] [1 + \epsilon_q], \quad (\text{A4})$$

where $E_e = q - Q_{\text{EC}} + m_e - \Delta_0$ and $\Delta_0 \simeq 90$ eV is a small correction introduced by Bahcall for the energy lost to

electronic rearrangement in the charge-changing reaction. The Gamow-Teller matrix element $B_{\text{GT}}^{(\nu,e)}(\text{gs})$ is determined from the measured half-life $\tau_{1/2}$ of the electron capture decay of ${}^{71}\text{Ge}$:

$$\omega = \frac{\ln[2]}{\tau_{1/2}} = \frac{G_F^2 \cos^2 \theta_C g_A^2}{2\pi} |\phi_{1s}|_{\text{avg}}^2 q_{1s}^2 \left[2(1 + \epsilon_o^{1s})(1 + \frac{P_L + P_M}{P_K}) \right] [2 B_{\text{GT}}^{(\nu,e)}(\text{gs})] [1 + g_{v,b(\text{EC})}] [1 + \epsilon_q]. \quad (\text{A5})$$

In this expression, $\epsilon_o^{1s} \simeq 0.013$ is a correction for overlap and exchange to account for the differences in the wave functions of the initial and final atomic states, $g_{v,b}$ is a radiative correction, and $\epsilon_q \simeq 5 \times 10^{-4}$ is a correction for

the weak magnetism contribution to the transition probability. The factor 2 with $B_{\text{GT}}^{(\nu,e)}(\text{gs})$ is a spin statistical factor. Hence,

$$\sigma_{\text{gs}} = \omega \frac{p_e E_e \mathcal{F}(Z_f, E_e)}{|\phi_{1s}|_{\text{avg}}^2 q_{1s}^2} [1 + g_{v,b(\text{EC})}]^{-1} \left[2(1 + \epsilon_o^{1s})(1 + \frac{P_L + P_M}{P_K}) \right]^{-1} \quad (\text{A6})$$

$$\equiv \sigma_0 \frac{p_e E_e \mathcal{F}(Z_f, E_e)}{m_e^2 c^3} \frac{1}{2\pi\alpha Z}. \quad (\text{A7})$$

The last line extracts the Bahcall cross-section scale σ_0 and inserts the dimensionful parameters needed to scale the energy and momentum. The atomic factors of Elliott *et al.* (2023) are equivalent to those of Bahcall:

$$\sum_i |\phi_i|_{\text{avg}}^2 q_i^2 = 0.01443 \frac{(m_e c^2)^5}{4\pi(\hbar c)^3} \quad (\text{A8})$$

Unfortunately, none of the three most advanced calculations referenced in (Bahcall, 1997) of the densities g_i has been published so the numerical factor 0.01443 is taken from the ratio in Eq. A3. (The computer code GRASP (Dyall *et al.*, 1989) may still be available.) Rather than use the theoretical electron density of each subshell at the nucleus, Elliott *et al.* use the experimentally measured subshell ratios, which has the advantage that only a single overlap and exchange correction and a single energy are needed, for the $1s$ state. As Bahcall has stated, overlap and exchange have little effect on the *total* EC rate, so Elliott *et al.* write Eq. A8 as follows, with the inclusion of a small Q-value update,

$$\sum_i |\phi_i|_{\text{avg}}^2 q_i^2 (1 + \epsilon_0^i) = |\phi_{1s}|_{\text{avg}}^2 q_{1s}^2 \left[(1 + \epsilon_o^{1s})(1 + \frac{P_L + P_M}{P_K}) \right] \quad (\text{A9})$$

The resulting expression is,

$$\sigma_0 = \frac{8\pi^2 \ln(2)\alpha Z}{|\phi_{1s}|_{\text{avg}}^2 q_{1s}^2 t_{1/2}} [1 + g_{v,b(\text{EC})}]^{-1} \left[2(1 + \epsilon_o^{1s})(1 + \frac{P_L + P_M}{P_K}) \right]^{-1}. \quad (\text{A10})$$

$$= \left(\frac{\hbar c}{m_e c^2} \right)^3 \frac{4\pi^2 \ln(2)\alpha Z}{0.01440 c t_{1/2}} \left(\frac{0.2221 \text{ MeV}}{q_{1s}} \right)^2 [1 + g_{v,b(\text{EC})}]^{-1} \quad (\text{A11})$$

$$= 8.63 \times 10^{-46} [1 + g_{v,b(\text{EC})}]^{-1} \text{ cm}^2, \quad (\text{A12})$$

which is only 0.3% larger than the value given in Bahcall (1997).

The above cross-section expressions give a ground-state cross section 2.6% lower than in the cross section table, Table II in Bahcall (1997). Whether this extends to the solar neutrino rates has not yet been evaluated, but in any scenario some reconsideration of the low-energy solar neutrino data may be required once a resolution to the Ga anomaly has emerged.

B. Bayesian Methods

The treatment of uncertainties when dealing with multiple data sets was addressed in the appendix of SF II, including the Inflation Factor Method used by the Particle Data Group that was adopted by most of the SF II working groups. Readers are directed there. We now extend that discussion to include the Bayesian methods used by several of the SF III work groups. For example,

in the analysis of the ${}^2\text{H}(p,\gamma){}^3\text{He}$ reaction presented in Section IV.D, both Bayesian analysis and Bayes model averaging were utilized.

Consider an optimization that utilizes Gaussian distributed priors in a constrained Bayesian analysis. With such priors, maximizing the likelihood amounts to minimizing an augmented χ^2

$$\chi_{\text{aug}}^2 \equiv \sum_{D,i} \frac{[y_{D_i} - f_D S(E; \lambda)]^2}{\sigma_{D_i}^2 + (\sigma_{D_i}^{\text{ext}})^2} + \sum_p \left(\frac{\lambda_p - \tilde{\mu}_p}{\tilde{\sigma}_p} \right)^2. \quad (\text{B1})$$

The first term in the above equation is the standard χ^2 and the second term derives from the Bayesian constraint on the parameters (λ_p) with prior mean ($\tilde{\mu}_p$) and width ($\tilde{\sigma}_p$), which are chosen with some prior knowledge or can be optimized as described below. Of course, care must be taken when choosing values for the priors ($\tilde{\mu}_p$ and $\tilde{\sigma}_p$). In the absence of prior information, it is common to set $\tilde{\mu}_p = 0$. Similarly, one should not make $\tilde{\sigma}_p$ too small, unless one has prior knowledge to do so, else this will bias the determination of λ_p .

In the first term in Eq. (B1), the double sum runs over the datasets (D) and individual results from each dataset (D_i), with the mean value and stochastic uncertainty of each data point given by y_{D_i} and σ_{D_i} , respectively.

The theoretical model describing the S-factor data, $S(E; \lambda)$, is a function of the energy (E) and a set of parameters (λ) that must be determined. The quoted systematic uncertainties are parameterized by the normalization factors f_D , with a prior of unit normalization and a width characterized by the quoted systematic uncertainty. It is straightforward within a Bayesian framework to utilize distribution functions for f_D that are not Gaussian, such as a log-normal or other distributions. The parameters f_D represent a normalization of the model function for a given dataset, which, viewed from a non-Bayesian perspective, can also be interpreted as normalization factors that must be applied to the data to match the “true” underlying distribution.

Finally, $\sigma_{D_i}^{\text{ext}}$ are unknown *extrinsic uncertainties* (de Souza *et al.*, 2019). If one assumes that a smooth function of energy can accurately describe the data, then the data in a given experimental set may scatter about this presumed “true” value by more than is reflected by the quoted statistical uncertainties. It was suggested in de Souza *et al.* (2019) that this extra scatter might be explained by some additional source of statistical uncertainty unbeknownst to the experimenter, or by some unknown systematic uncertainty that is different for each data point in the same data set, as opposed to a correlated systematic that affects all data points similarly. In either case, this extrinsic uncertainty can be accommodated in a Bayesian analysis framework by adding the additional uncertainty as a normal-distributed source of noise with a width that is constrained by the data. In de Souza *et al.* (2019) and Moscoso *et al.* (2021), $\sigma_{D_i}^{\text{ext}}$

was added as an absolute uncertainty, independent of the energy. For some data, Odell *et al.* (2022a) suggested adopting instead a relative uncertainty, such that the scale of the extrinsic fluctuations is proportional to the mean value of $S(E)$. This strategy of adding extra extrinsic uncertainty to the data sets can be viewed as an alternative to that of inflating the quoted statistical uncertainties by $\sqrt{\chi_\nu^2}$, which is often used for seemingly incompatible data sets (Workman *et al.*, 2022), where χ_ν^2 is the χ^2 per degree of freedom. An advantage of the extrinsic uncertainty method is that it uses the observed scatter within a given data set as a measure of the possible size of unreported uncertainties, as opposed to uniformly increasing the uncertainty in all data sets by the same relative amount.

One can further perform a Bayes model averaging, which we implement and describe here. For a given model, after optimizing the posterior parameter distributions, the Bayes factor (BF) is proportional to the probability of the model given the data (Hilbe *et al.*, 2017). Therefore, for a fixed data set, the BF can be used as a relative probability of each model, thus enabling a weighted model-averaging procedure. If we assume a uniform likelihood for each model, the expectation value and variance of a quantity Y is given by

$$\text{E}[Y] = \sum_k \text{E}[Y|M_k]P[M_k|D] \quad (\text{B2})$$

$$\begin{aligned} \text{Var}[Y] &= \sum_k \text{Var}[Y|M_k]P[M_k|D] \\ &\quad + \sum_k \text{E}^2[Y|M_k]P[M_k|D] - \text{E}^2[Y], \end{aligned} \quad (\text{B3})$$

where $\text{E}[Y|M_k]$ denotes the expectation of Y given the model M_k and $P[M_k|D]$ denotes the probability of the model M_k given the data (D), which is given by

$$P[M_k|D] = \frac{\text{BF}_{M_k}}{\sum_l \text{BF}_{M_l}}, \quad (\text{B4})$$

with BF_{M_l} the Bayes factor of model l . Similarly, for a given model, in the absence of prior information on the size of an unknown parameter, the optimal width of its prior can be estimated by finding the value of $\tilde{\sigma}_p$ that maximizes the BF, which typically provides a reasonable approximation to marginalizing over the prior width.

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Solar Fusion III: supplementary material

Data and analysis for the astrophysical S-factor of the ${}^7\text{Be}(p, \gamma){}^8\text{B}$ reaction

We present additional information related to the Solar Fusion III section on the astrophysical S-factor of the ${}^7\text{Be}(p, \gamma){}^8\text{B}$ reaction, S_{17} .

I. Experimental Data

In this section we reproduce the data that formed the basis for the S_{17} recommendation of the Solar Fusion III evaluation. We provide the original data in Tables I, II, III, IV, V, VI, VII, VIII, IX, X, and XI. The common mode errors associated with each data set are shown in Table XII. For the Coulomb dissociation datasets we provide the estimated asymmetric common mode errors, inflated to account for the uncertainty in the $E2$ components as described in section B of the main text.

TABLE I S_{17} data of Baby *et al.* (2003)^a 5% added as per a private communication reported in SF II

E_{cm} (keV)	S_{17} (eV b)	Error (eV b)
302	18.1	2 ^a
356	18.8	1.4 ^a
415	20.2	1.8 ^a
844	23.6	0.8
849	23.8	0.8
853	23.8	0.6
856	24.3	0.6
1078	25.5	0.8

TABLE II S_{17} data of Buompane *et al.* (2022)

E_{cm} (keV)	S_{17} (eV b)	Error (eV b)
367.2	14.7	2.43
610.5	51.2	8
632.2	86.1	9
632.2	87.2	10.5
632.5	90.9	19
667.2	25.9	5.1
697.8	20	7.4
699.8	30.7	9
812.2	25.3	5.1

II. Effective Field Theory EFT_{gs} and EFT* fits

Here we provide more details on some of the fits that were performed using EFT_{gs} and EFT* in the $S_{17}(0)$ extrapolations. We performed fits in three different energy regions with and without Coulomb dissociation (CD) data. For notational purposes, we define fits to radiative capture data for $E \leq 475$ keV region I, for $E \leq 1250$ keV

TABLE III S_{17} data of Filippone *et al.* (1983)

E_{cm} (keV)	S_{17} (eV b)	Error (eV b)
117	18.9	2.6
186	17.4	2.1
257	18.4	2
302	16.7	1.6
346	19.2	1.5
390	20.9	1.3
435	19.5	1.5
480	17.4	1
524	20.4	1.1
568	25.2	1.4
590	33.5	1.7
603	43.3	2.3
612	59.4	2.8
625	89.3	4.4
632	101.5	2.8
638	93.6	4.1
648	65.5	3.3
656	52.8	2.5
678	30.3	1.6
701	27.9	1.4
745	22.6	1.1
789	24.8	1.6
877	21.3	1.4
1053	20.5	1.1
1230	23.8	1.4

TABLE IV S_{17} data of Hammache *et al.* (1998)

* Data points with $E > 1250$ keV are not included in any fit.

E_{cm} (keV)	S_{17} (eV b)	Error (eV b)
353.8	16.7	1.6
364.1	17.9	1.7
403.7	20	1.8
451.2	17.7	1.6
496	17.9	1.8
881.2	20.6	1.7
899.2	21.3	2.1
899.3	20.9	2
1096.7	19.2	1.7
1397.3	23.1	2
500.3	17.1	1.6
880.5	19	1.9
1199.7	20.5	2.1
1310.4*	20.6	2.1

region II, and the non-resonant energy range comprised of $E \leq 490$ keV and 805 keV $\leq E \leq 1250$ keV region III. When CD data is included in the fits in the three energy regions we denote them as IB, IIB and IIIB, respectively. Since EFT_{gs} is only physically relevant below the ${}^7\text{Be}^*$ excitation energy $E_* = 429.1$ keV, it is used only in the I and IB fits but provides valuable checks for the EFT* II

TABLE V S_{17} data of Hammache *et al.* (2001)

E_{cm} (keV)	S_{17} (eV b)	Error (eV b)
111.7	15.8	2.7
134.7	19.5	3.1
185.8	17.2	2.1

TABLE VI S_{17} data of Hass (1999)

* Data points with $E > 1250$ keV are not included in any fit.

E_{cm} (keV)	S_{17} (eV b)	Error (eV b)
1090.0	22.7	1.2
1290.0*	23.8	1.5

and IIB fits. Frequentist χ^2 minimization and Bayesian fits were performed at LO, NLO and NNLO. We will only report the NNLO numbers to keep the amount of information manageable.

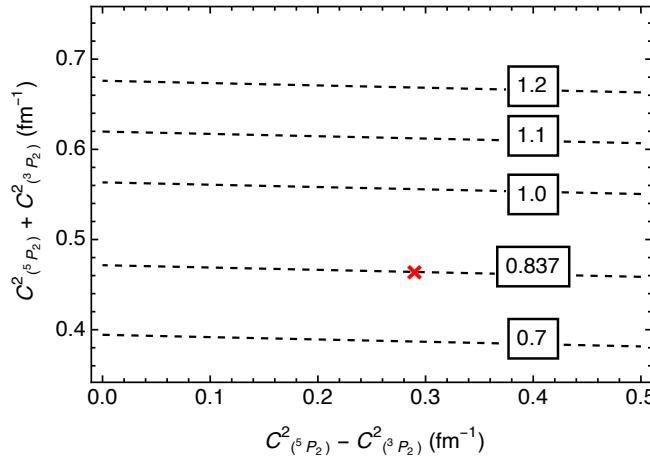


FIG. 1 Contour plot of $S_{17}(0)$ in EFT_{gs} (Higa *et al.*, 2022) normalized to 21 eV b. The boxed numbers identify the dashed contour lines. The \times marks the experimental determinations of the ANC (Tabacaru *et al.*, 2006; Trache *et al.*, 2003).

The extrapolations of S_{17} to 50 eV from the various fits of radiative capture data using χ^2 minimization are shown in Table XIII. We include a $\sqrt{2/\text{dof}}$ uncertainty estimate in the reduced χ^2 values. All the χ^2 fits included a theory prior for the ANCs in the 3P_2 and ${}^3P_2^*$ channels and a prior on the s -wave scattering length ratio a_{22}/a_{12} as described in the main text. This ensured that the fitted parameter values are consistent with the EFT power counting. Fits without the theory priors, indicated as I^a, work just as well. However, the fitted parameters, shown in Table XIV, can take on values that would invalidate the EFT power counting and expansion. For the EFT_{gs} I^a fit, this is not a major concern. This theory has only two fit parameters and fits without priors give values close to expectations; see Table XIV. For the EFT_{*} I^a

TABLE VII S_{17} data of Junghans *et al.* (2010)

* Data points with $E > 1250$ keV are not included in any fit.

E_{cm} (keV)	S_{17} (eV b)	Error (eV b)
116.1	20.2	1.1
140.1	18.9	0.8
184.3	19.7	0.7
255.5	19.4	0.5
277.8	19.8	0.5
326.7	19.7	0.4
362.2	20.7	0.5
872.7	24.7	0.6
184.6	19.9	0.5
220	19.4	0.5
255.6	19.6	0.4
277.7	20.2	0.4
326.6	20.8	0.4
362.2	20.3	0.3
871.5	24.6	0.3
999.5	24.9	0.4
1099.9	25.9	0.3
1200.1	26.7	0.6
1754.2*	31.1	0.8

TABLE VIII S_{17} data of Strieder *et al.* (2001)

* Data points with $E > 1250$ keV are not included in any fit.

E_{cm} (keV)	S_{17} (eV b)	Error (eV b)
322	17.8	1.4
364	16.8	1
461	15.7	0.9
548	23.4	1.5
771	20.9	1.1
859	20.4	1.8
903	20.4	1.8
991	20	1
1122	22	1.9
1254	19.2	1
1386	21.6	1.1
1736*	25.2	1.2
2610*	36.7	2.3

that has 4 fit parameters, the uncertainties in the fitted parameters are very large. Moreover, the central values result in a negligible contribution from the 3P_2 channel with a much larger contribution from the ${}^3P_2^*$ channel achieved through an unnaturally large ANC and a_{22}/a_{12} ratio. Despite the large uncertainties in the parameters, the $S_{17}(0)$ values are similar to the fits constrained by theory priors.

We can understand the robustness of the $S_{17}(0)$ values and their small errors despite the large parameter uncertainties in the I^a fit using Fig. 1. The nearly horizontal contours show that the S-factor at threshold is mostly sensitive to the sum of the squares of the ANCs in EFT_{gs}. The capture cross sections in the 5P_2 and 3P_2 channels have the same momentum dependence. The different s -wave scattering length in the two channels affects the momentum dependence, but only at NNLO. Thus the

TABLE IX S_{17} Data of Schümann *et al.* (2006)

E_{cm} (keV)	S_{17} (eV b)	Error (eV b)
160	17.5	2.1
316	19.3	1.2
942	23.6	1.6
1244	25.2	1.9

TABLE X S_{17} Data of Davids and Typel (2003)

E_{cm} (keV)	S_{17} (eV b)	Error (eV b)
192	15.2	1.0
384	16.1	0.9

$E1$ transition can be equally described through capture in either of the two spin channels leading to a single parameter family of theories constrained only by the total sum of the squares of the ANCs $C_{(5P_2)}^2 + C_{(3P_2)}^2 = C_{p_{3/2}}^2 + C_{p_{1/2}}^2$. In EFT*, the mixing in the $S = 1$ channels makes the analysis more complicated. However, we still find correlations among the 5P_2 , 3P_2 and ${}^3P_2^*$ channels such that $C_{(5P_2)}^2 + C_{(3P_2)}^2$ is well constrained, as shown in the last column of Table XIV.

Table XV shows the data set scaling factors for EFT* I^a and I fits that had very different values for the fitted parameters. Again, we see that there is no difference in the quality of the fits. The difference is in how the fits without theoretical constraints enhance contributions from the ${}^3P_2^*$ channel at the expense of the 3P_2 channel. This imbalance is slightly less pronounced when we fit over a wider energy range where the contributions from the excited ${}^7\text{Be}^*$ become more important, better constraining the scattering length ratio a_{22}/a_{12} . However, sensitivity to the initial choice of parameter values in the χ^2 minimization remains since in the EFT, a_{22}/a_{12} is still a NNLO effect even at the higher energies.

In Table XVI, we show the $S_{17}(0)$ extrapolations that include the CD data in the χ^2 minimization at NNLO. All the fits were with theory priors. We find that including CD data results in a slightly lower $S_{17}(0)$ that is in agreement with the radiative capture fits in Table XIII. The scaling factors for the IIB fit are shown in Table XV.

We emphasize that in the EFT framework it is not sufficient that the theory fit experimental data accurately, but must do so in a manner that is consistent with the power counting such that the perturbative expansion remains meaningful. Without a consistent expansion, theoretical uncertainty estimates are not reliable. This leads us to consider Bayesian estimations of $S_{17}(0)$.

The χ^2 minimizations demonstrate that both EFT_{gs} and EFT* are able to describe data self-consistently in their respective domain of validity, once the theory priors are included. Incorporating EFT priors in the fits is natural in the Bayesian estimations.

TABLE XI S_{17} Data of Kikuchi *et al.* (1998)

E_{cm} (keV)	S_{17} (eV b)	Error (eV b)
365	18.0	0.4
865	21.1	1.0
1118	23.6	1.2

TABLE XII Common mode errors of each experiment.

^a 2.3% common mode error plus 0.37% (the largest of the BE3L and BE3S data sets) varying error from Table I of Jung-hans *et al.* (2010)

DATA	Common Mode Error (%)
Baby <i>et al.</i> (2003)	2.2
Buompane <i>et al.</i> (2022)	4
Filippone <i>et al.</i> (1983)	11.9
Hammache <i>et al.</i> (1998)	5.7
Hammache <i>et al.</i> (2001)	11.5
Hass (1999)	5
Junghans <i>et al.</i> (2010)	2.67 ^a
Strieder <i>et al.</i> (2001)	8.3
Schümann <i>et al.</i> (2006) (GSI)	(-7.7, +5.6)
Davids and Typel (2003) (MSU)	(-7.1, +9.6)
Kikuchi <i>et al.</i> (1998) (RIKEN)	(-12.7, +8.4)

Moreover, CD data is found to be compatible with the radiative capture data. The scaling factors in the fits are also reasonable. Therefore, the central consideration was whether one should fit data in only region IB, as was done in SF II without CD data, using the simpler EFT_{gs} with only two fit parameters, or one should use data over the larger range of region IIB with the more complete EFT*.

The $S_{17}(0)$ estimation from Bayesian analysis is shown in Table XVII. The EFT_{gs} and EFT* fits in the low energy region IB are nearly identical. The Bayesian “evidence”—the theory posterior probability distribution given some data D and hypothesis H —comparison gives $\ln[P(\text{EFT}_{\text{gs}}|D, H)/P(\text{EFT}_*|D, H)] \approx 0.32(33)$, ever so slightly favoring the simpler 2-parameter theory. However, this is not strong evidence given the systematic uncertainty of around 2 in the log-evidence calculations (Higa *et al.*, 2022). We also calculate the average

TABLE XIII $S_{17}(50 \text{ eV})$ and its first two energy derivatives at NNLO (Higa *et al.*, 2022) determined from χ^2 minimization with only radiative capture data. The theoretical uncertainty is not included.

Theory	S_{17} (eV b)	$\frac{S'_{17}}{S_{17}}$ (MeV $^{-1}$)	$\frac{S''_{17}}{S_{17}}$ (MeV $^{-2}$)	χ^2
EFT _{gs} I ^a	20.3(5)	-1.79(5)	31.3(7)	0.96(25)
EFT _{gs} I	20.3(5)	-1.79(5)	31.3(7)	0.95(25)
EFT* I ^a	20.8(8)	-1.85(27)	31.3(17)	1.02(26)
EFT* I	20.8(7)	-1.84(11)	31.3(10)	0.93(25)
EFT* II	20.7(4)	-1.86(4)	31.3(6)	1.08(16)
EFT* III	20.8(4)	-1.84(4)	31.3(6)	1.08(18)

TABLE XIV The mean and standard deviation of EFT (Higa *et al.*, 2022) parameters from χ^2 minimization. For comparison, the expected sizes of parameters from the ANCs are: $C_{(5P_2)}^2 = 0.376(16) \text{ fm}^{-1}$, $C_{(3P_2)}^2 = 0.088(7) \text{ fm}^{-1}$ and $C_{(3P_2^*)}^2 = 0.1215(36) \text{ fm}^{-1}$ (Tabacaru *et al.*, 2006; Trache *et al.*, 2003; Zhang *et al.*, 2018). $\bar{L}_{22} \sim 1$ is evaluated at $p = p_R$. The last column is not a fit parameter but just lists the sum of ANC squares $C^2 = C_{(5P_2)}^2 + C_{(3P_2)}^2$. Fits are only to radiative capture data.

Theory	$C_{(5P_2)}^2 (\text{fm}^{-1})$	$C_{(3P_2)}^2 (\text{fm}^{-1})$	$C_{(3P_2^*)}^2 (\text{fm}^{-1})$	a_{22}/a_{12}	$ \bar{L}_{22} $	$r_s (\text{MeV})$	$C^2 (\text{fm}^{-1})$
EFT _{gs} I ^a	0.421 ± 0.032	0.128 ± 0.046	—	—	—	—	0.549(20)
EFT _{gs} I	0.444 ± 0.015	0.093 ± 0.017	—	—	—	—	0.537(13)
EFT _* I ^a	0.552 ± 1.620	0 ± 2	0 ± 10^6	0 ± 10^7	—	—	0.552(28)
EFT _* I	0.463 ± 0.027	0.088 ± 0.018	0.121 ± 0.013	-0.701 ± 0.833	—	—	0.551(20)
EFT _* II	0.460 ± 0.018	0.089 ± 0.017	0.121 ± 0.013	-0.971 ± 0.276	2.55(10)	$-110.9(1)$	0.549(10)
EFT _* III	0.470 ± 0.017	0.082 ± 0.017	0.121 ± 0.013	-0.833 ± 0.267	—	—	0.551(10)

TABLE XV Scaling factors for data sets at NNLO (Higa *et al.*, 2022) from χ^2 minimization.

Data set	EFT _* I ^a	EFT _* I	EFT _* IIB
Baby	1.00(2)	1.00(2)	0.97(2)
Buompane	1.02(4)	1.02(4)	1.07(3)
Filippone	1.01(4)	1.01(4)	1.03(2)
Hammache '98	1.05(4)	1.05(4)	1.10(3)
Hammache '01	1.05(8)	1.05(8)	1.06(4)
Hass	—	—	1.02(3)
Junghans BE3	0.95(2)	0.95(2)	0.93(1)
Strieder	1.15(5)	1.15(4)	1.12(3)
GSI	—	—	0.99(4)
MSU	—	—	1.15(5)
RIKEN	—	—	1.05(2)

TABLE XVI $S_{17}(50 \text{ eV})$ and its first two energy derivatives determined from χ^2 minimization (Higa *et al.*, 2022) including CD data. The theoretical uncertainty is not included.

Theory	S_{17} (eV b)	$\frac{S'_{17}}{S_{17}}$ (MeV^{-1})	$\frac{S''_{17}}{S_{17}}$ (MeV^{-2})	χ^2
EFT _{gs} IB	20.1(4)	-1.79(5)	31.3(7)	0.95(23)
EFT _* IB	20.5(7)	-1.84(11)	31.3(10)	0.93(23)
EFT _* IIB	20.5(3)	-1.86(3)	31.3(5)	1.03(15)
EFT _* IIIB	20.8(4)	-1.84(3)	31.3(5)	0.96(17)

$\langle \chi^2 \rangle$ from the posterior distribution to signify qualitatively the goodness of fit for the Bayesian estimates since the evidence cannot be used to compare the quality of fits when the data sets in IB and IIB are different. The standard deviation in the χ^2 posterior is negligible and so we only quote a $\sqrt{2/\text{dof}}$ uncertainty for $\langle \chi^2 \rangle$.

In Fig. 2 we compare the three Bayesian fits over a wide energy range. The uncertainties in the fits are very small (see Table XX) and are not shown. The $M1$ contribution from the 1^+ resonance is only included in the IIB fit that covers the resonance region. The Bayesian estimates of the EFT parameters are in Table XVIII. The parameter values are consistent with the EFT power counting. We find correlations among the 5P_2 , 3P_2 , and ${}^3P_2^*$ channels

TABLE XVII $S_{17}(50 \text{ eV})$ and its first two energy derivatives at NNLO determined from Bayesian analysis (Higa *et al.*, 2022) including CD data. The theoretical uncertainty is not included.

Theory	S_{17} (eV b)	$\frac{S'_{17}}{S_{17}}$ (MeV^{-1})	$\frac{S''_{17}}{S_{17}}$ (MeV^{-2})	$\langle \chi^2 \rangle$
EFT _{gs} IB	20.1(4)	-1.79(5)	31.9(8)	1.13(28)
EFT _* IB	20.5(5)	-1.84(10)	31.9(12)	1.24(29)
EFT _* IIB	20.5(3)	-1.86(4)	31.9(6)	1.06(16)

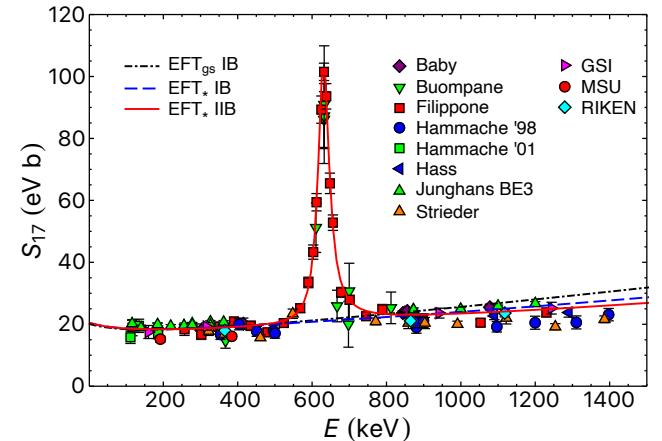


FIG. 2 Astrophysical S-factor for ${}^7\text{Be}(p, \gamma){}^8\text{B}$ at NNLO from Bayesian fits. The dot-dashed (black) curve is the $E1$ contribution from the EFT_{gs} IB fit, dashed (blue) curve is the $E1$ contribution from the EFT* IB fit and the solid (red) curve is from the EFT* IIB fit that includes both the $E1$ and $M1$ contributions (Higa *et al.*, 2022).

such that only $C_{(5P_2)}^2 + C_{(3P_2)}^2$ is well constrained, as in the χ^2 minimization fits.

The posterior from the EFT* IIB fit is shown in Fig. 3 as an example. The correlations between the ANCs support the general picture shown in Fig. 1. The interpretation in EFT_{gs} is simpler than in EFT* where the mixing in the $S = 1$ channels forces a correlation with the s -wave

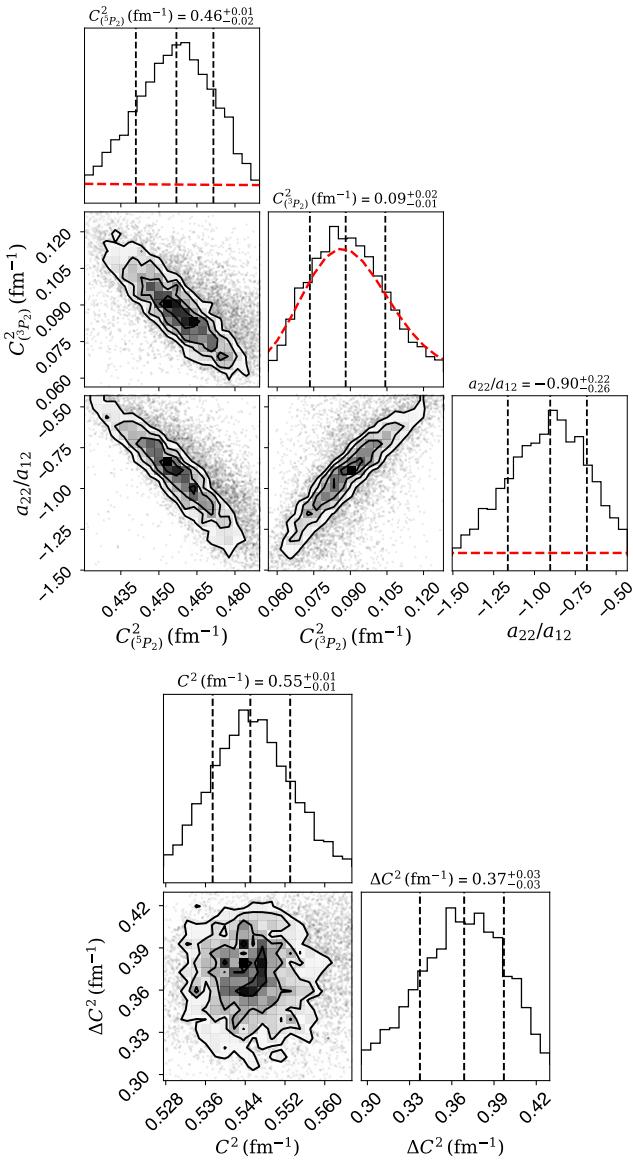


FIG. 3 Posterior distributions for the EFT* IIB fit at NNLO (Higa *et al.*, 2022). In the top panel, the dashed (red) curves are the priors. In the bottom panel, $C^2 = C_{(5P_2)}^2 + C_{(3P_2)}^2$ and $\Delta C^2 = C_{(5P_2)}^2 - C_{(3P_2)}^2$. 99% of the posterior probability distribution is displayed. The vertical lines indicate the median and the regions containing 68% of the posterior probability around it.

scattering length ratio a_{22}/a_{12} as well. Nonetheless, the sum of ANC squares $C^2 = C_{(5P_2)}^2 + C_{(3P_2)}^2$ is well constrained in Fig. 3 and Table XVIII. Based on the parameter estimates in Table XVIII, one might conclude that the fits determine the individual ANCs well. However, this is not accurate.

In the top panel of Fig. 3 we see that the $C_{(3P_2)}^2$ posterior is nearly identical to its prior. Something similar happens for $C_{(3P_2^*)}^2$ as well. On the other hand, we do gain

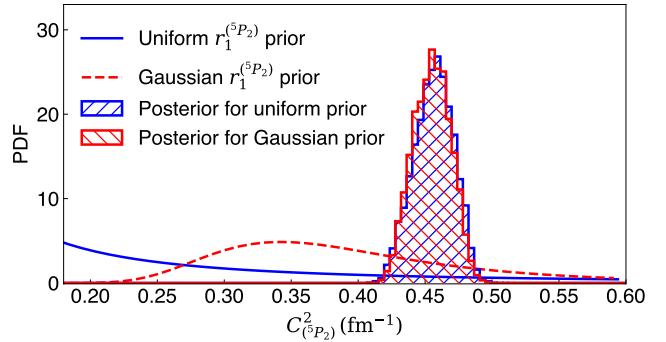


FIG. 4 Probability distribution functions (PDFs): Two different priors and the corresponding posteriors for $C_{(5P_2)}^2$ (Higa *et al.*, 2022). Details are in the text.

information about $C_{(5P_2)}^2$ from the fits, as seen from the corresponding prior and posterior. In EFT_{gs} and EFT*, the 5P_2 channel is dominant, thus, the information from the capture which we expect to be sensitive only to the sum $C^2 = C_{(5P_2)}^2 + C_{(3P_2)}^2$ is channeled into the posterior of $C_{(5P_2)}^2$, while keeping the $C_{(3P_2)}^2$ and $C_{(3P_2^*)}^2$ posteriors fixed near their priors. Similar observations apply to the χ^2 fits with theory priors in Table XIV.

In Fig. 4 we compare two $C_{(5P_2)}^2$ posteriors. One was obtained from a uniform prior on the p -wave effective range $r_1^{(5P_2)} \sim \mathcal{U}(-100 \text{ MeV}, 1.5 \text{ MeV})$ that covers the physically relevant regions. The $C_{(5P_2)}^2$ prior was obtained for the plot from the $r_1^{(5P_2)}$ prior using:

$$P_Y(y) = P_X[g(y)]g'(y), \quad (1)$$

where $P_X(x)$, $P_Y(y)$ are the probability distribution functions for the random variables X and Y , respectively, and $g(y) = f^{-1}(y)$ for $y = f(x)$, an increasing function of x . Uniform priors on EFT parameters are a reasonable choice, as often we only estimate the sizes of the couplings and not their signs. Here, the sign of the ANC squares sets an upper bound on $r_1^{(5P_2)}$ (Higa *et al.*, 2022). A wide Gaussian prior $r_1^{(5P_2)} \sim \mathcal{N}(-40 \text{ MeV}, 10 \text{ MeV})$ centered around the experimental ANC is also shown. The Gaussian prior can result in an unphysical $r_1^{(5P_2)} \gtrsim 1.5 \text{ MeV}$ but with a insignificant cumulative probability that is less than 2×10^{-5} . The priors result in nearly identical $C_{(5P_2)}^2$ posteriors, confirming the robustness of the fits.

The data set scaling factors are shown in Table XIX. We find the three fits to be consistent with each other. The scaling factors from χ^2 minimization that includes the priors in the cost function are similar to the ones from the Bayesian estimates as seen, for example, for the IIB fits in Tables XV and XIX. Further, in Table XX we include the $E1$ and $M1$ contributions from the EFT* IIB Bayesian fit at NNLO at various energies. We find the

$S_{17}(E)$ posterior pdf to be close to a Gaussian, and so only list the mean and standard deviations.

III. A Second Bayesian Halo EFT Analysis

Next we provide some details about the set of fits that apply the EFT theory and Bayesian analysis tool developed in (Zhang *et al.*, 2015, 2018) to analyze the data sets summarized in Sec. I. This EFT (Zhang *et al.*, 2015) describes non-resonant capture up to NLO in terms of an EFT expansion. It has 9 parameters, including two ANCs of the final bound state: $C_{(3P_2)}$ and $C_{(5P_2)}$ for the 3P_2 and 5P_2 channels, a parameter ϵ_1 describing the relative strength of the radiative capture to the excited ${}^7\text{Be}$ core configuration in the final bound state, two s -wave scattering length parameters: $a_{(3S_1)}$ and $a_{(5S_2)}$, two s -wave effective range parameters: $r_{(3S_1)}$ and $r_{(5S_2)}$, and two parameters describing two-body-current-type contributions in the $S = 1$ and 2 channels: \bar{L}_1 and \bar{L}_2 .

Zhang *et al.* (2018) demonstrates that this NLO EFT can represent previous model calculations with a few percent accuracy at energies up to 1 MeV. This few-percent discrepancy can serve as a simple estimate of the theoretical errors of this NLO EFT, assuming that (1) the previous models, when represented in the EFT framework, can populate beyond-NLO terms in the EFT expansion, and (2) neglecting them in the NLO EFT is the source of the observed discrepancies.

To describe the resonant component in the capture data, we add a Breit-Wigner (BW) parameterization (Rolsf and Rodney, 1988) to the EFT-based non-resonant S-factor. The BW parameterization has three parameters: resonance energy E_{res} , the proton decay width $\Gamma_p(E_{\text{res}})$, and the radiative width $\Gamma_\gamma(E_{\text{res}})$, and takes the following form:

$$S_{\text{res}}(E) = \frac{3\pi}{8k^2} \frac{\Gamma_p(E)\Gamma_\gamma(E)E e^{2\pi\eta(E)}}{(E - E_{\text{res}})^2 + (\Gamma_{\text{tot}}(E)/2)^2}, \quad (2)$$

$$\Gamma_p(E) = \Gamma_p(E_{\text{res}}) \frac{e^{-2\pi\eta(E)}}{e^{-2\pi\eta(E_{\text{res}})}}, \quad (3)$$

$$\Gamma_\gamma(E) = \Gamma_\gamma(E_{\text{res}}) \frac{E_\gamma^3}{E_{\gamma\text{res}}^3}, \quad (4)$$

where $\eta(E)$ is the Sommerfeld parameter and E_γ the energy of the emitted photon at energy E .

In some fits, the frequentist approach of χ^2 minimization was explored. We found that some of the fitted results depended on the starting point in the model parameter space for the χ^2 -minimization algorithm; the computation of the Hessian matrix needed for error propagation in the S-factor calculations could become ill-conditioned, which renders numerical calculation unstable. The Bayesian analysis, on the other hand, encountered no such difficulties. Therefore, we report only on the Bayesian analysis. Three different choices of energy

ranges for the direct capture measurements have been explored: region I with $E \leq 0.49$ MeV, II with $E \leq 1.25$ MeV, and III with $E \leq 0.49$ MeV and $0.805 \leq E \leq 1.25$ MeV. We then added the Coulomb-disassociation (CD) data sets in our analysis to the capture data sets with I, II, and III energy ranges. They are labeled as IB, IIB, and IIIB, following the convention in Sec. II. We note that in the I, III, IB, and IIIB fits, our model only includes non-resonant EFT components, while for the II and IIB fits, we also include a BW resonance.

It is an essential advantage that Bayesian analysis allows the transparent incorporation of assumptions or previous knowledge about the physical model by specifying the prior probability distributions for model parameters. For this EFT, we choose wide ranges, as allowed by existing information. We assume $C_{(3P_2)}^2$ and $C_{(5P_2)}^2$ to have uniform prior distributions between 0 and 1 fm^{-1} ; $r_{(3S_1)}$ and $r_{(5S_2)}$ to have uniform prior distributions between 0 and 10 fm ; ϵ_1 to have a uniform distribution between -1 and 1 ; \bar{L}_1 and \bar{L}_2 to have uniform distributions between -10 and 10 fm ; $a_{(3S_1)}$ and $a_{(5S_2)}$ to have Gaussian distributions centered on the recently experimentally inferred central values of 17.34 and -3.18 fm respectively (Paneru *et al.*, 2019), with standard deviations equal to the experimental errors of 1.33 and 0.55 fm , respectively (Paneru *et al.*, 2019). We further enforce constraints on s -wave scattering parameters that preclude the existence of s -wave resonances below 0.6 MeV , following the same treatment as Zhang *et al.* (2015). For the resonance parameters, we assume uniform distributions between 0.6 and 0.7 MeV for E_{res} , 0.003 and 0.063 MeV for $\Gamma_p(E_{\text{res}})$, and 1×10^{-10} and $4 \times 10^{-8}\text{ MeV}$ for $\Gamma_\gamma(E_{\text{res}})$. These represent 5σ ranges (Buompane *et al.*, 2022).

As mentioned in the main text, the other parameters to be fitted are the scaling factors for each experiment. Their prior distributions are Gaussian distributions centered at 1 with standard deviations equal to the common-mode errors (CMEs) listed in Table XII. The prior distributions of the scaling factors of the CD experiments are asymmetric Gaussians because their CMEs are asymmetric.

Table XXI summarizes the median and 1σ error for the predicted S-factor and its derivatives at $E = 0\text{ MeV}$ from six different fits. All these fits show consistency and stability. For each experiment, the fitted mean value and 1σ errors for the scaling factor minus 1 are listed in Table XXII in percent for the I, II, and III analyses and in Table XXIII for the analysis with CD data included. A negative mean value suggests a preference for scaling the data downward, while a positive mean suggests upward scaling.

The mean and error for the EFT parameters in all the fits are listed in Table XXIV. As can be seen, the results are consistent and stable against changes to the fitted energy range and inclusion of CD data. Table XXV

TABLE XVIII The median and the interval containing 68% of the posterior probability of the EFT (Higa *et al.*, 2022) parameters at NNLO from Bayesian fits. CD data are included in the fits. The last column lists $C^2 = C_{(5P_2)}^2 + C_{(3P_2)}^2$.

Theory	$C_{(5P_2)}^2(\text{fm}^{-1})$	$C_{(3P_2)}^2(\text{fm}^{-1})$	$C_{(3P_2^*)}^2(\text{fm}^{-1})$	a_{22}/a_{12}	$ \bar{L}_{22} $	$r_s(\text{MeV})$	$C^2(\text{fm}^{-1})$
EFT _{gs} IB	$0.438^{+0.012}_{-0.013}$	$0.0932^{+0.0146}_{-0.0151}$	—	—	—	—	$0.531^{+0.01}_{-0.009}$
EFT _* IB	$0.477^{+0.025}_{-0.024}$	$0.0866^{+0.0156}_{-0.0145}$	$0.12^{+0.011}_{-0.011}$	$-0.56^{+0.622}_{-0.74}$	—	—	$0.542^{+0.017}_{-0.015}$
EFT _* IIB	$0.479^{+0.016}_{-0.017}$	$0.0881^{+0.0162}_{-0.0147}$	$0.12^{+0.012}_{-0.011}$	$-0.904^{+0.225}_{-0.26}$	$3.03^{+0.07}_{-0.06}$	$-110.9^{+0.2}_{-0.1}$	$0.545^{+0.008}_{-0.008}$

TABLE XIX Bayesian estimates of scaling factors for the data sets at NNLO (Higa *et al.*, 2022).

Data set	EFT _{gs} IB	EFT _* IB	EFT _* IIB
Baby	1.00(2)	1.00(2)	0.97(1)
Buompane	1.02(3)	1.02(4)	1.08(3)
Filippone	1.00(3)	1.00(3)	1.03(2)
Hammache '98	1.05(3)	1.05(3)	1.11(3)
Hammache '01	1.06(4)	1.07(4)	1.06(4)
Hass	1.00(3)	1.00(3)	1.02(3)
Junghans BE3	0.94(2)	0.94(2)	0.93(1)
Strieder	1.15(4)	1.15(4)	1.13(2)
GSI	0.99(4)	0.99(4)	0.99(3)
MSU	1.16(4)	1.17(4)	1.16(4)
RIKEN	1.06(3)	1.06(3)	1.05(2)

shows the mean and error of $C_{(3P_2)}^2 + C_{(5P_2)}^2$, which is the key quantity determining $S(0)$. As we can see, including the CD data sets reduces this quantity somewhat and thus reduces $S(0)$, but these results are consistent with those excluding the CD data sets. It should be noted that individual ANCs are only weakly constrained, as shown in Fig. 5, but the two are strongly anti-correlated such that the sum of their squares is tightly constrained.

The means and errors of the resonance parameters are provided in Table XXVI. The energy of the resonance is tightly constrained, while the widths are constrained on the percent level. Again, the two fits are consistent.

Fig. 6 compares the IIB fit and the experimental data sets used in the analysis. The energy range shown exceeds that of the fit. Some of the errors are smaller than the size of the symbols. The theoretical 1σ error band is only visible near $E = 1.2$ MeV. It should be pointed out that the experimental data points have not been renormalized by the scaling factors s_α . The $S(E)$ curves derived from the other 5 fits are similar and are not shown here.

IV. R-matrix Theory Fits

We investigated the fractional contribution of the 2.2 MeV, 3^+ resonance to the S-factor using R-matrix calculations. As illustrated in Fig. 7, it is vanishingly small in the energy range of interest.

In the rest of this section, we attempt to quantify the uncertainty of the R-matrix fit of the low-energy

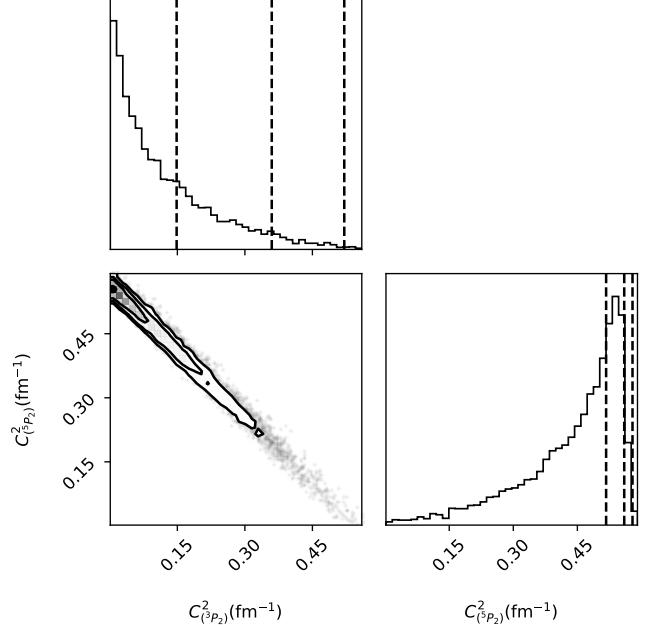


FIG. 5 Correlation plot of $C_{(3P_2)}^2$ and $C_{(5P_2)}^2$ from the Zhang *et al.* (2015) EFT + BW fit of the IIB experimental data. The contours enclose 68%, 95%, and 99.7% of the posterior probability from the inside out. In the 1-dimensional projections, the corresponding confidence intervals all start from 0 and end at one of the dashed lines. $C_{(3P_2)}^2$ and $C_{(5P_2)}^2$ are strongly anti-correlated such that their sum is tightly constrained (see Table XXV).

$^7\text{Be}(p, \gamma)^8\text{B}$ capture data. The fit includes all experiments that have data below 425 keV and the parameters correspond to the two ANCs and four background pole partial widths. The normalization of each dataset is allowed to float following the prescription described in the main text. We randomly select $N = 2300$ starting points in this parameter space and for each we perform a χ^2 minimization. As the problem exhibits multiple local minima, most runs finish the fit procedure at a different one. Therefore, in what follows the usual definition for

TABLE XX The $E1$ and $M1$ contributions to S_{17} from the Bayesian fit of EFT_{*} to data in region IIB at NNLO (Higa *et al.*, 2022) including CD data. The means and standard deviations are listed. An additional 3% NNLO theoretical uncertainty applies and should be added in quadrature.

E (keV)	$E1\ S_{17}$ (eV b)	$M1\ S_{17}$ (eV b)
0.05	20.54(28)	0.000 184 7(61)
3.224	20.42(28)	0.000 207 6(68)
11.27	20.15(28)	0.000 267 3(86)
24.19	19.78(27)	0.000 381 2(120)
41.98	19.38(26)	0.000 587 3(183)
64.65	19.01(26)	0.000 958 1(296)
92.20	18.71(25)	0.001 632(50)
124.6	18.49(25)	0.002 874(88)
161.9	18.38(24)	0.005 215(160)
204.1	18.36(24)	0.009 755(299)
251.1	18.44(24)	0.018 924(580)
303.1	18.61(24)	0.038 578(1184)
359.9	18.86(24)	0.084 71(261)
421.5	19.19(24)	0.2103(65)
488.1	19.57(25)	0.6586(207)
496.7	19.62(25)	0.7812(247)
505.5	19.67(25)	0.9349(297)
514.3	19.73(25)	1.131(36)
523.2	19.78(25)	1.384(45)
532.1	19.83(25)	1.718(56)
541.2	19.89(25)	2.168(71)
550.3	19.94(25)	2.792(94)
559.5	20.00(25)	3.684(126)
568.8	20.05(25)	5.012(177)
578.1	20.11(25)	7.084(261)
587.5	20.17(25)	10.51(41)
597.0	20.22(25)	16.6(7)
606.6	20.28(25)	28.17(129)
616.3	20.34(26)	50.22(237)
626.0	20.39(26)	79.51(291)
635.8	20.45(26)	79.06(260)
645.7	20.51(26)	51.66(215)
655.6	20.57(26)	31.23(130)
665.7	20.63(26)	20.00(79)
675.8	20.69(26)	13.76(51)
686.0	20.75(26)	10.05(36)
696.2	20.81(26)	7.693(263)
706.6	20.87(26)	6.111(204)
717.0	20.93(26)	4.999(163)
727.5	21.00(26)	4.187(135)
738.0	21.06(26)	3.577(114)
748.7	21.12(26)	3.105(98)
759.4	21.19(26)	2.733(86)
770.2	21.25(27)	2.435(76)
781.1	21.32(27)	2.191(68)
792.0	21.38(27)	1.989(61)
803.0	21.45(27)	1.819(56)
814.1	21.52(27)	1.676(51)
825.3	21.58(27)	1.553(47)
836.5	21.65(27)	1.448(44)
847.9	21.72(27)	1.356(41)
859.3	21.79(27)	1.276(39)
870.8	21.86(28)	1.205(36)
882.3	21.93(28)	1.143(34)
893.9	22.01(28)	1.088(33)
989.7	22.62(29)	0.8033(239)
1090	23.29(31)	0.6658(196)
1196	24.03(33)	0.5927(173)
1306	24.84(36)	0.5533(160)

Energy Range	$S(0)$ (eV b)	$S'/S(\text{MeV}^{-1})$	$S''/S(\text{MeV}^{-2})$
I	21.02(75)	-1.832(119)	31.90(29)
II	21.18(50)	-1.919(69)	32.03(19)
III	21.14(51)	-1.862(58)	32.01(12)
IB	20.71(71)	-1.825(112)	31.89(25)
IIB	21.03(50)	-1.915(71)	32.02(19)
IIIB	20.96(49)	-1.861(59)	32.01(12)

TABLE XXI The S-factor and its 1st and 2nd derivatives at $E = 0$ from 6 different fits using the Zhang *et al.* (2015) EFT.

Data sets	I	II	III
Junghans BE3	-3.1(1.9)	-3.7(1.7)	-2.6(1.7)
Filippone	3.9(3.4)	5.2(2.1)	8.2(2.5)
Hammache '98	4.1(3.5)	10.4(2.4)	10.6(2.5)
Baby	-0.18(2.1)	-2.4(1.7)	-2.1(1.7)
Boumpane	2.2(3.8)	8.4(3.1)	2.0(3.8)
Strieder	13.2(3.4)	12.3(2.2)	14.8(2.4)
Hass	NA	2.8(3.7)	4.1(3.7)

TABLE XXII Fitted values of $100 \times (s_\alpha - 1)$, with s_α the scaling factor for the data set α , from the Bayesian analysis of radiative capture data in energy ranges I, II, and III using the Zhang *et al.* (2015) EFT.

the mean and standard deviation are replaced by

$$\langle x \rangle = \sum_{i=1}^N w_i x_i / \sum_{i=1}^N w_i$$

$$\sigma = \sqrt{\sum_{i=1}^N w_i (x_i - \langle x \rangle)^2 / \sum_{i=1}^N w_i}, \quad (5)$$

where the weight of each sample is calculated based on the respective χ^2 value: $w_i = \exp[-(\chi^2/\chi_{\min}^2)^2]$. The value χ_{\min}^2 is the minimum χ^2 obtained from all the samples and appears there only to normalize the weights. By squaring the exponent we achieve a flatter, broader Gaussian, so as to avoid assigning an overwhelmingly large weight to the best fitting calculation. The resulting squared ANCs from all the calculations are shown in Figure 8, with the plotted lines showing the mean and standard deviation of $C_1^2 + C_2^2$ computed using Eq. (5). It is clear that what is constrained by the fit presented here is the sum of the squares of the ANCs and not each of them separately, but it is worth noting that in the majority of fits, $C_2 > C_1$. The overall result is $S(0) = 19.8 \pm 0.8$ eV b, consistent with our recommendation.

The ANCs determine the overall normalization of the S-factor but, since we are allowing for floating normalizations of the data sets, the overall normalization we are trying to reproduce is also somewhat varying. For this set of calculations the value of χ^2 , without dividing by the number of degrees of freedom, clusters around a value of 7, with a minimum at 6.4, but calculations with val-

Data sets	IB	IIB	IIIB
Junghans BE3	-3.9(1.9)	-4.3(1.6)	-3.5(1.7)
Filippone	3.2(3.4)	4.7(2.1)	7.7(2.4)
Hammache '98	3.7(3.5)	10.1(2.4)	10.2(2.6)
Baby	0.02(2.1)	-2.6(1.6)	-2.7(1.6)
Boumpane	2.0(3.9)	8.2(3.3)	2.0(3.7)
Strieder	12.5(3.5)	11.9(2.2)	14.3(2.4)
Hass	NA	2.8(3.6)	4.1(3.6)
GSI	-0.3(4.4)	-0.62(3.5)	0.62(3.38)
MSU	15.3(3.5)	15.4(3.6)	16.1(3.5)
RIKEN	5.4(2.7)	5.7(2.2)	6.9(2.2)

TABLE XXIII Fitted values of $100 \times (s_\alpha - 1)$, with s_α the scaling factor for the data set α , from the Bayesian analysis of radiative capture and Coulomb dissociation data in energy ranges I, II, and III using the [Zhang et al. \(2015\)](#) EFT.

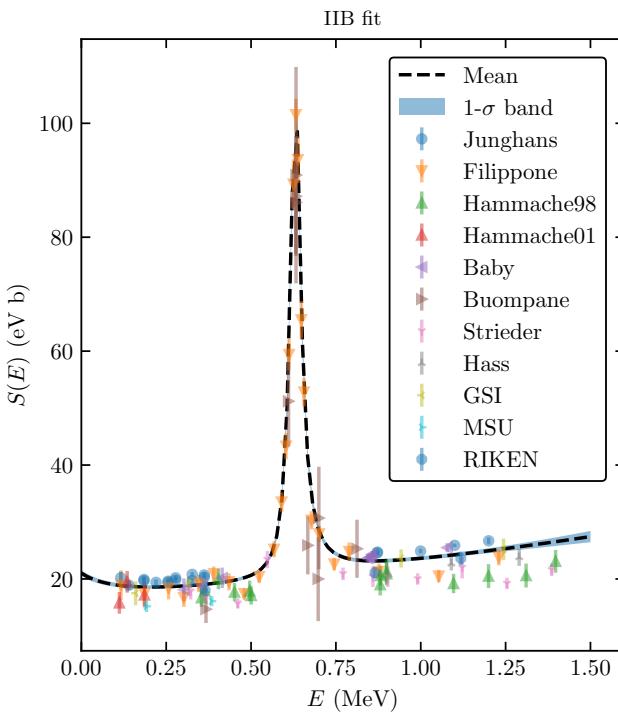


FIG. 6 The [Zhang et al. \(2015\)](#) EFT + BW fit of the IIB experimental data, including both radiative capture and CD.

ues exceeding 20 exist. We find that most calculations are concentrated around the same value of the ANC (given by the slope in Figure 8) but lie on the rightmost edge of the uncertainty of the value extracted from ${}^8\text{B}$ breakup using both a Glauber model description ([Trache et al., 2001](#)) and continuum-discretized coupled channel calculations ([Sürer et al., 2022](#)), as shown in the blue histogram in Figure 9. In a stark shift, the probability is concentrated at larger values of the sum of the squared ANCs and exhibits a significantly smaller spread after the ANC values obtained through the R-matrix fit are

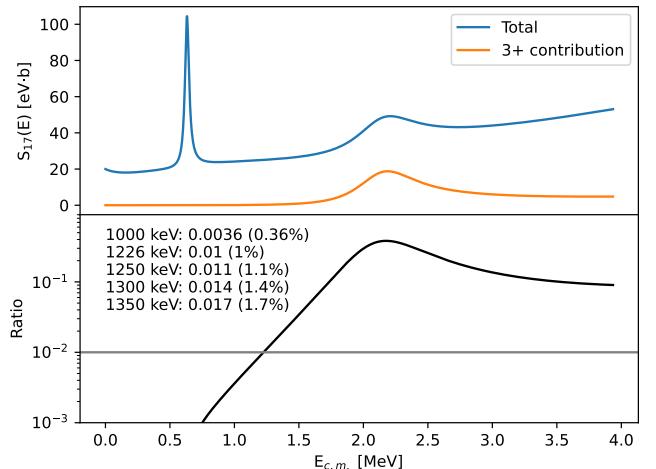


FIG. 7 The contribution of the 2.2 MeV, 3^+ resonance and the total of all contributions to the astrophysical S-factor $S_{17}(E)$ of a representative R-matrix calculation are shown in both absolute and relative terms. The 3^+ contribution is negligibly small in the energy range of our fits.

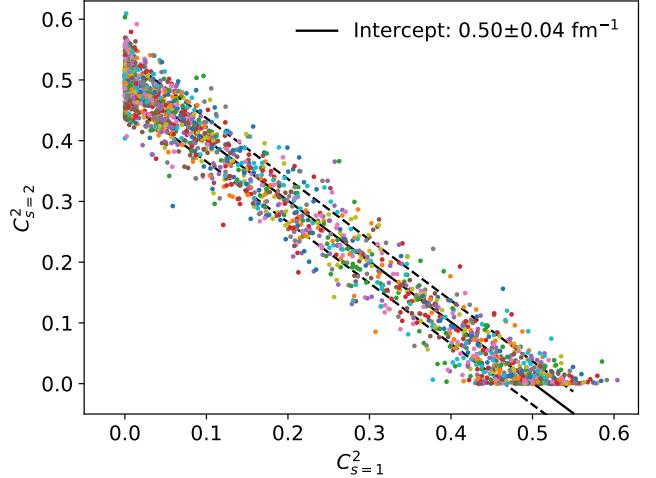


FIG. 8 Squared ANCs in fm^{-1} of each of the N calculations after performing a χ^2 minimization to determine the R-matrix fit parameters.

divided by the normalization of the Junghans data set, as can be seen in the red histogram of Figure 9. This result graphically depicts the tension between the normalization of the Junghans data set and the analyses of [Trache et al. \(2001\)](#) and [Sürer et al. \(2022\)](#).

It should be noted that a similar strong covariance exists between the ANCs and many other data sets, for example, dividing by the RIKEN data set normalization enhances the agreement with the breakup analyses and dividing by the Baby data leaves the histograms mostly unchanged. From this set of fits we can obtain the weighted mean and standard deviation for the the first and second

Analysis	$C_{(3P_2)}^2(\text{fm}^{-1})$	$a_{(3S_1)}$ (fm)	$r_{(3S_1)}$ (fm)	ϵ_1	\bar{L}_1 (fm)	$C_{(5P_2)}^2(\text{fm}^{-1})$	$a_{(5S_2)}$ (fm)	$r_{(5S_2)}$ (fm)	\bar{L}_2 (fm)
I	0.22(15)	17.2(1.4)	3.8(2.4)	0.1(6)	1.6(3.6)	0.34(16)	-3.2(6)	5.1(2.9)	1.5(5.4)
II	0.12(11)	17.2(1.3)	2.4(1.6)	0.2(6)	2.0(3.8)	0.45(11)	-3.2(5)	5.1(2.8)	6.7(2.4)
III	0.07(8)	17.2(1.3)	3.6(2.2)	0.0(6)	2.8(3.6)	0.49(9)	-3.2(6)	3.9(2.9)	5.4(2.1)
IB	0.21(15)	17.2(1.3)	3.9(2.4)	0.1(6)	1.8(3.6)	0.34(15)	-3.2(5)	5.2(2.9)	1.7(5.3)
IIB	0.12(11)	17.2(1.3)	2.4(1.6)	0.2(6)	2.0(3.8)	0.44(12)	-3.2(6)	5.2(2.8)	6.6(2.5)
IIIB	0.07(8)	17.3(1.3)	3.6(2.2)	0.0(6)	2.8(3.7)	0.48(9)	-3.2(6)	3.8(2.8)	5.3(2.3)

TABLE XXIV The mean values and 1σ errors of the NLO EFT parameters from all six fits using the [Zhang et al. \(2015\)](#) EFT.

analysis	$C_{(3P_2)}^2 + C_{(5P_2)}^2(\text{fm}^{-1})$	$S(0)(\text{eV}\cdot\text{b})$
I	0.556(25)	21.02(75)
II	0.565(16)	21.18(50)
III	0.561(16)	21.14(51)
IB	0.548(23)	20.71(71)
IIB	0.561(16)	21.03(50)
IIIB	0.556(15)	20.96(49)

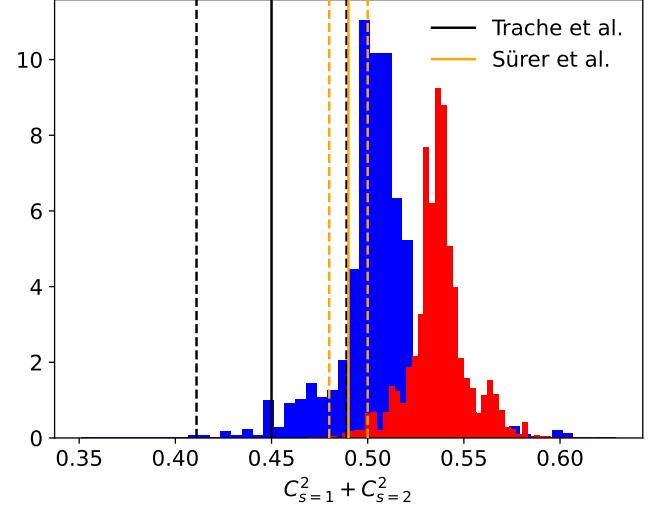
TABLE XXV The mean values and 1σ errors of the sum of the squares of the ANCs from all six fits using the [Zhang et al. \(2015\)](#) EFT. The $S(0)$ values are also listed.

Analysis	$E_{\text{res}}(\text{MeV})$	Γ_p (MeV)	Γ_γ (MeV)
II	0.6315(6)	$3.86(15) \times 10^{-2}$	$2.73(12) \times 10^{-8}$
IIB	0.6315(6)	$3.86(15) \times 10^{-2}$	$2.72(12) \times 10^{-8}$

TABLE XXVI Fitted values of the 3 BW resonance parameters in the II and IIB analyses, in which the [\(Zhang et al., 2015\)](#) EFT is used to fit the direct capture component.

derivatives of the S-factor at zero energy $S'(0)/S(0) = -1.8 \pm 0.1 \text{ MeV}^{-1}$ and $S''(0)/S(0) = 39.1 \pm 0.7 \text{ MeV}^{-2}$. The value of the first derivative is in agreement with that given in the main text, which is based on more data taken over a wider energy range, while the second derivative lies well outside the corresponding 3σ range.

The combined set of all the S-factor curves calculated using R-matrix fits is shown in Figure 10 as black lines with a transparency inversely proportional to their weight; the weighted mean $\pm 1\sigma$ is also shown in red. From the fairly constant size of the standard deviation of the mean at various energies, we can postulate a high degree of correlation between the values of the S-factor at different energies. This is indeed the case, with correlation coefficients across all energies exceeding 0.9. The implication here is, in turn, that a precise measurement, even at high energies, will have a significant impact on the evaluation as a whole. This impact is also demonstrated in Figure 10, with a precise, fictitious data point at 0.35 MeV lying on top of the evaluated mean and having an uncertainty of 0.3 eV·b. Recomputing the weights associated with each fit we obtain a mean (blue) that is almost on top of the previous one but with significantly

FIG. 9 Posterior distribution of the sum of the squares of the computed ANC values with (blue) and without (red) renormalization of the Junghans data set in the R-matrix analysis. The solid and dashed lines give the mean and 1σ uncertainty of the sum of the squares of the ANC values extracted from ${}^8\text{B}$ breakup by [Trache et al. \(2001\)](#) and [Suerer et al. \(2022\)](#).

reduced uncertainty. This high degree of correlation is a direct byproduct of the small number of parameters that completely determine the capture cross section. Extending the analysis to the full energy spectrum (up to 1.25 MeV) would introduce more parameters (for example to describe the $M1$ resonance) and thus would reduce the correlation somewhat, but the main argument presented here would still stand. It would therefore be highly desirable for future evaluations to benefit from such experiments, should they become available, and for theoretical approaches to seek an even more accurate description of the process at experimentally accessible energies.

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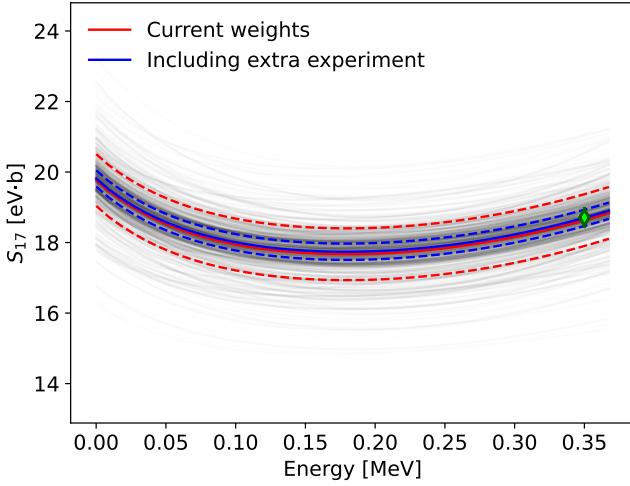


FIG. 10 S-factor realizations obtained from different starting points in the R-matrix fit (see text for more details). The red solid line represents the mean of the evaluation with the 1σ band denoted by dashed lines. Introducing a precise, fictitious experimental point at 0.35 MeV (in green) significantly reduces the evaluation uncertainty (blue).

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