

Cosmological variation of the deuteron binding energy, strong interaction, and quark masses from big bang nucleosynthesis

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We use big bang nucleosynthesis calculations and light element abundance data to constrain the relative variation of the deuteron binding energy since the Universe was a few minutes old, $\delta Q = Q(BBN) - Q(\text{present})$. Two approaches are used, first treating the baryon to photon ratio η as a free parameter, but with the additional freedom of varying δQ , and second using the WMAP value of η and solving only for δQ . Including varying Q yields a better fit to the observational data than imposing the present day value, rectifying the discrepancy between the ${}^4\text{He}$ abundance and the deuterium and ${}^7\text{Li}$ abundances, and yields good agreement with the independently determined η_{WMAP} . Using η_{WMAP} , the minimal deviation consistent with the data is significant at about the 4σ level; $\delta Q/Q = -0.019 \pm 0.005$. If the primordial ${}^4\text{He}$ abundance lies towards the low end of values in the literature, this deviation is even larger and more statistically significant. Taking the light element abundance data at face value, our result may be interpreted as variation of the dimensionless ratio $X = m_s/\Lambda_{\text{QCD}}$ of the strange quark mass and strong scale: $\delta X/X = (1.1 \pm 0.3) \times 10^{-3}$. These results provide a strong motivation for a more thorough exploration of the potential systematic errors in the light element abundance data.

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I. INTRODUCTION

Recent astronomical data suggest a possible variation of the fine structure constant $\alpha = e^2/\hbar c$ at the 10^{-5} level over a time scale of 10 billion years, see Ref. [1] (a discussion of other limits can be found in Ref. [2], and references therein). Naturally, these data motivated more general discussions of possible variations of other constants. Unlike for the electroweak forces for the strong interaction, there is generally no direct relation between the coupling constants and observable quantities. In recent papers [3–5], we presented general discussions on the possible influence of the strong scale variation on primordial big bang nucleosynthesis (BBN) yields, the Oklo natural nuclear reactor, quasar absorption spectra and atomic clocks. Here we continue this work, concentrating on BBN.

One can only measure variations of dimensionless parameters [2]. Big bang nucleosynthesis is sensitive to a number of fundamental dimensionless parameters including the fine structure constant α , $\Lambda_{\text{QCD}}/M_{\text{Plank}}$, and m_q/Λ_{QCD} , where m_q is the quark mass and Λ_{QCD} is the strong scale determined by a position of the pole in the perturbative QCD running coupling constant. In this work we search for any possible variation of m_q/Λ_{QCD} because there is a mechanism which provides a very strong sensitivity of BBN to this parameter.

The first and most crucial step in BBN is the process $p + n \rightarrow d + \gamma$. The synthesis starts at $t \geq 3$ sec when the temperature goes down below $T \leq 0.6$ MeV and lasts until t

≤ 6 min when the temperature becomes $T \leq 0.05$ MeV. The reaction rate for the above process defines all subsequent processes and final primordial abundances of light elements. Amongst the factors that can influence the reaction rate, the most significant seems to be a variation of the deuteron binding energy (this variation was discussed in Refs. [3,5–12]). Indeed, the equilibrium concentration of deuterons and the inverse reaction rate depend exponentially on it. Moreover, the deuteron is a shallow bound level. Therefore the relative variation of the deuteron binding Q is much larger than the relative variation of the strong potential U , i.e., $\delta Q/Q \gg \delta U/U$. As a result the variations in the strong interaction may be most pronounced via the deuteron binding energy. We also take into account the effect of variation of the virtual level in the neutron-proton system [13], which is even more sensitive to the variation of the strong interaction.

The question we address here is whether or not existing observations of the primordial abundances of the light elements suggest any change in the deuteron binding energy at the time of BBN. To do so, we use a compilation of light element abundance data from the literature for ${}^4\text{He}$, ${}^7\text{Li}/\text{H}$ and D/H . As we show later, the currently greater experimental precision on ${}^4\text{He}$ results in that element dominating our results. The other 2 light elements nevertheless provide important consistency checks.

The data we use for ${}^4\text{He}$ is presented in Table I and comprised 14 surveys giving estimates for the primordial value Y_p , the mass fraction, derived using, or by extrapolation to,

TABLE I. Data on the primordial ${}^4\text{He}$ mass fraction.

Y_p	Ref.
0.2391 ± 0.0020	[14]
0.2384 ± 0.0025	[15]
0.2371 ± 0.0015	[16]
0.2443 ± 0.0015	[17]
0.2351 ± 0.0022	[18]
0.2345 ± 0.0026	[19]
0.244 ± 0.002	[20]
0.243 ± 0.003	[21]
0.232 ± 0.003	[22]
0.240 ± 0.005	[23]
0.234 ± 0.002	[24]
0.244 ± 0.002	[25]
0.242 ± 0.009	[26]
0.2421 ± 0.0021	[27]

low metallicity in each case. There is clear evidence for significant scatter amongst these 14 values, presumably due to unquantified systematics, or if not, intrinsic inhomogeneities. The dominance by ${}^4\text{He}$, or indeed by any single element, unfortunately increases susceptibility to systematic errors, and we have therefore attempted to explore the effect of these in several ways.

First, in order to make best use of all the available ${}^4\text{He}$ data, we add a constant term to each of the statistical errors on Y_p , such that the normalized χ^2 for all 14 points about the weighted mean value is equal to unity. This approach is equivalent to the assumption that all 14 estimates of Y_p are unbiased and Gaussian distributed, but that there is an additional systematic component to the statistical error which is different (and hence random) for each estimate.

Second, as shown later, smaller values of Y_p are less consistent with $\delta Q/Q=0$ than larger values. Thus we carry out a reanalysis using a subset of the Y_p 's, taking only the highest values such that the normalized χ^2 about the weighted mean value is equal to unity, *without* increasing the individual errors by a constant, as described above. This procedure selects 9 values from the original 14. In doing this, we are exploring the consequence of there being strong systematics for the small Y_p 's, and little or none for the high values. This is conservative, in the sense that we are minimizing our estimate for $\delta Q/Q$.

Finally, in order to obtain some estimate of the plausible range on our estimate of $\delta Q/Q$, we perform the converse analysis, subsetting the data by discarding *high* values of Y_p , again such that the normalized χ^2 about the weighted mean value is equal to unity. This leaves 9 points. The two samples thus overlap.

The data on deuterium abundances D/H from quasar absorption systems were selected according to two criteria:

(i) Metallicity must be low, so as to more closely reflect primordial values [Si/H] or [O/H] less than or equal to -2.0 .

(ii) Must be detection, not upper limit.

TABLE II. Data on the primordial deuterium abundance.

QSO	$z(\text{abs})$	$D/H \times 10^{-5}$	[Si/H]	Ref.
Q1009+299	2.504	4.0 ± 0.65	-2.53	[28]
PKS1937-1009	3.572	3.25 ± 0.3	-2.26 [O/H]	[29]
HS0105+1619	2.536	2.5 ± 0.25	-2.0	[30]
Q2206-0199	2.076	1.65 ± 0.35	-2.23	[31]
Q1243+3047	2.526	$2.42 + 0.35 - 0.25$	-2.77 [O/H]	[32]

These requirements leave only five data points listed in Table II.

The data for lithium primordial abundance are shown in Table III. Here $A = \log(Y_{\text{Li}}) + 12$, where $Y_{\text{Li}} = {}^7\text{Li}/\text{H}$. Applying the first procedure described above, in order to obtain $\chi^2/N=1$, we have to add to the individual σ 's 0.0017 for helium points, 0.344×10^{-5} for deuterium points, and 0.028 for lithium points. For the weighted mean values we obtain

$$Y_p = 0.2393 \pm 0.0011, \quad (1)$$

$$Y_D = D/H = (2.63 \pm 0.31) \times 10^{-5}, \quad (2)$$

and

$$A = 2.315 \pm 0.051. \quad (3)$$

The latter value corresponds to the following lithium abundance:

$$Y_{\text{Li}} = (2.02 \pm 0.22) \times 10^{-10}. \quad (4)$$

The second and the third procedures are meaningful only for the helium points. The number of deuterium points is too small and the lithium data points are the least scattered. We need only 20% increase in individual uncertainties to bring χ^2/N to 1 for the lithium data. In addition, the deuterium and the lithium data do not produce a significant contribution in determination of $\delta Q/Q$ which is entirely dominated by the helium data due to their high accuracy.

Keeping 9 upper points for the helium mass fraction data, that give $\chi^2/N=0.94$, we obtain for the weighted mean value

$$Y_p = 0.2424 \pm 0.0008. \quad (5)$$

If we keep 9 lower points, we obtain

TABLE III. Data on the primordial Li/H abundance.

A	Ref.
$2.09 + 0.11 - 0.12$	[33]
2.35 ± 0.1	[34]
2.36 ± 0.12	[35]
$2.34 \pm 0.056 \pm 0.06$	[36]
$2.07 + 0.16 - 0.04$	[37]
2.22 ± 0.20	[38]
2.4 ± 0.2	[39]
2.5 ± 0.1	[40]

$$Y_p = 0.2363 \pm 0.0008, \quad (6)$$

which is significantly lower than both in Eqs. (5) and (1).

II. THE BBN EQUATIONS

We use the standard BBN set of equations that describe the time development of the abundances of the elements in an expanding Universe [41]

$$\frac{\dot{R}}{R} = H = \sqrt{\frac{8\pi}{3M_p^2} \rho_T}, \quad (7)$$

$$\frac{\dot{n}_B}{n_B} = -3H, \quad (8)$$

$$\dot{\rho}_T = -3H(\rho_T + p_T), \quad (9)$$

$$\dot{Y}_i = \sum_{j,k,l} N_i \left(\Gamma_{kl \rightarrow ij} \frac{Y_l^{N_l} Y_k^{N_k}}{N_l! N_k!} - \Gamma_{ij \rightarrow kl} \frac{Y_i^{N_i} Y_j^{N_j}}{N_i! N_j!} \right), \quad (10)$$

$$n_- - n_+ = n_B \sum_j Z_j Y_j, \quad (11)$$

where n_B is the density of baryons and Y_i is the abundance of the element ${}^A Z_i$. The right-hand side of Eq. (10) corresponds to a reaction

$$N_i ({}^A Z_i) + N_j ({}^A Z_j) \leftrightarrow N_k ({}^A Z_k) + N_l ({}^A Z_l). \quad (12)$$

ρ_T and p_T denote total energy density and pressure, respectively,

$$\rho_T = \rho_\gamma + \rho_e + \rho_\nu + \rho_B, \quad (13)$$

$$p_T = p_\gamma + p_e + p_\nu + p_B. \quad (14)$$

Equation (7) defines the expansion rate. Equation (8) defines the change in time of the baryon density, and the rate equation (10) defines the time evolution of the abundances and their final values after freeze-out. Equation (11), where n_- and n_+ are the densities of electrons and positrons, is the condition of electroneutrality that defines a chemical potential of electrons.

III. EFFECT OF THE DEUTERON BINDING ENERGY VARIATION

The sensitivity of the reaction rates $\Gamma_{\gamma d \leftrightarrow pn}$ to parameters of the strong interaction in general, and to the deuteron binding energy in particular, comes from two sources. First, the reaction rate $\Gamma_{\gamma d \rightarrow pn}$ depends exponentially on the deuteron binding energy Q . Second, the cross section of the reaction $np \rightarrow \gamma d$ is very sensitive to the position of the virtual level with the energy $\epsilon_v = 0.07$ MeV. Any change in the strong NN potential causing a shift in the deuteron binding energy Q will change the position of the virtual level ϵ_v as well. The relation between δQ and $\delta \epsilon_v$ can be obtained using the fact

that both a real level and a virtual one are close to $E=0$. The relation is (see the Appendix)

$$\frac{\delta \epsilon_v}{\sqrt{\epsilon_v}} = - \frac{\delta Q}{\sqrt{Q}}. \quad (15)$$

The cross section for the $n + p \rightarrow d + \gamma$ reaction can be found in textbooks [13]. In the leading order in Q/ϵ_v the product of the cross section and the velocity is proportional to

$$\sigma v \sim Q^{5/2} / \epsilon_v.$$

Thus, in linear order in δQ we have the following modification of the reaction rate:

$$\Gamma_{np \rightarrow d\gamma} \rightarrow \Gamma_{np \rightarrow d\gamma} \left[1 + \left(5/2 + \sqrt{\frac{Q}{\epsilon_v}} \right) \frac{\delta Q}{Q} \right]. \quad (16)$$

We should note, however, that according to our BBN calculations the effect of the deuteron binding energy variation for the inverse reaction $d\gamma \rightarrow np$ is more important due to exponential dependence on Q of the reaction rate $\Gamma_{d\gamma \rightarrow np}$ than the variation of the cross section $np \rightarrow d\gamma$ given by Eq. (16). We have taken both effects into account anyway.

We modified one of the standard BBN codes [42] in such a way that Q can be changed for this reaction. Varying Q changes the abundances of all three elements under discussion. In Fig. 1 we plot the abundance of D, the mass fraction of ${}^4\text{He}$, and the abundance of ${}^7\text{Li}$, as functions of Q at the value of the baryon to photon ratio $\eta = 6.14 \times 10^{-10}$ found from anisotropy of cosmic microwave background [43]. From Fig. 1 we see that the deuterium abundance is not very sensitive to Q . The data are fully compatible with the present value of the deuteron binding energy. Such a poor sensitivity can be explained by relatively large error bars for the deuterium abundance.

The data on ${}^4\text{He}$, in contrast, show strong sensitivity to the deuteron binding energy favors lower Q during primordial nucleosynthesis. The data on ${}^7\text{Li}$ also favors lower Q approximately for the same δQ as ${}^4\text{He}$.

Figure 1 gives a qualitative picture of the dependence of light element abundances on the deuteron binding energy. In order to obtain more quantitative results we analyze the likelihood functions as functions of Q and η .

A. The likelihood functions

The likelihood function for the abundances have been chosen in the form

$$L_f(\eta, Q) = \exp \left(-\frac{1}{2} \sum_{ij} [Y_i^{\text{th}}(\eta, Q) - Y_i^{\text{ex}}] \times w_{ij} [Y_j^{\text{th}}(\eta, Q) - Y_j^{\text{ex}}] \right). \quad (17)$$

Here the sum goes over three light elements, w_{ij} is the inverse error matrix that was calculated using the approach

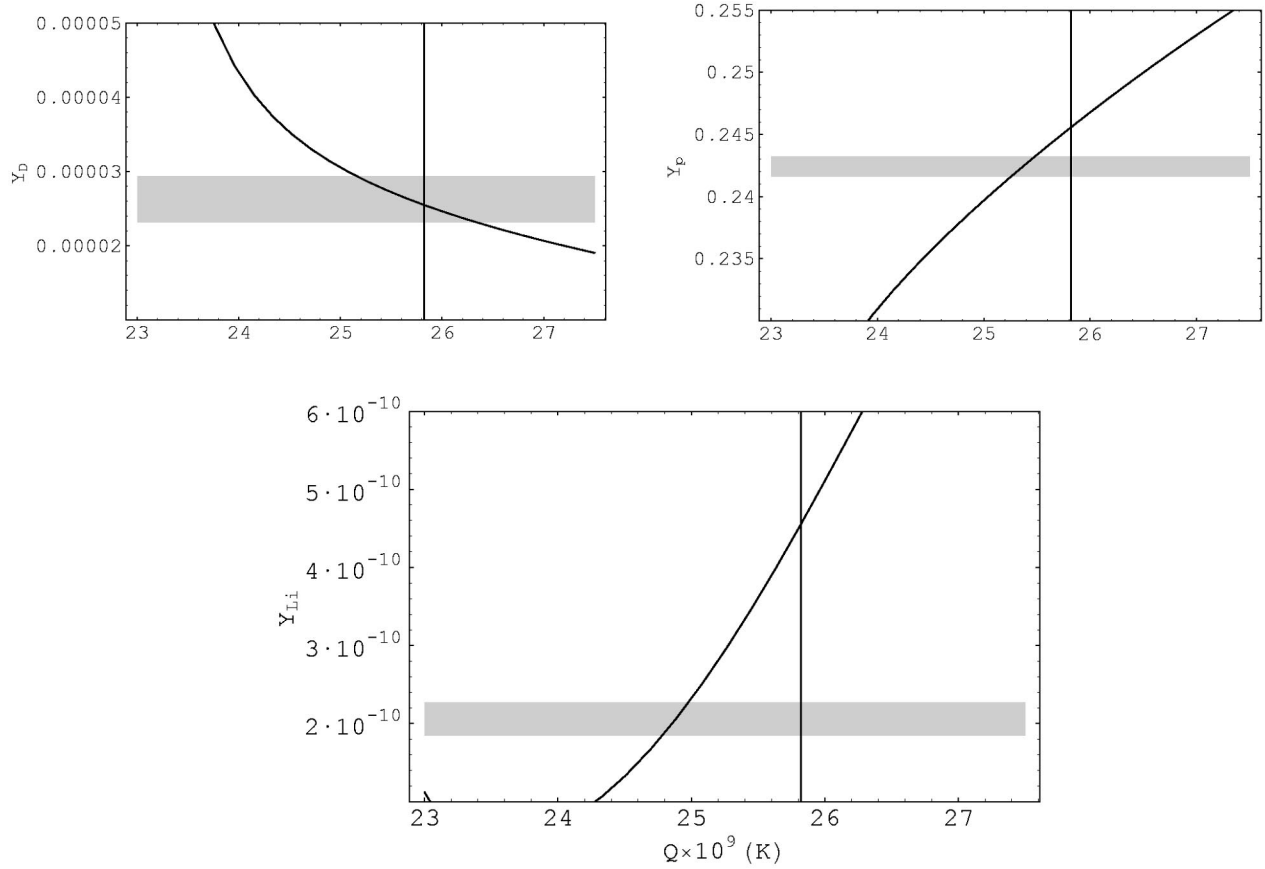


FIG. 1. The light elements abundances act as a function of the deuteron binding energy Q for $\eta_{\text{WMAP}} = 6.14 \times 10^{-10}$. The vertical line shows the present value of $Q = 25.82 \times 10^9$ K. The shaded regions illustrate the 1σ range in the data. For helium the high Y_p value [Eq. (5)] is shown.

proposed in Ref. [44]. The errors in theoretical values of the abundances can be found from the uncertainties in the reaction rates

$$\delta Y_i^{\text{th}} = Y_i^{\text{th}} \sum_k \lambda_{ik} \frac{\Delta R_k}{R_k}, \quad (18)$$

where ΔR_k are the reaction rate errors and

$$\lambda_{ik} = \frac{\partial \ln Y_i^{\text{th}}}{\partial \ln R_k}$$

are the logarithmic derivatives. The error matrix σ_{ij} can be calculated then by

$$\sigma_{ij}^2 = Y_i^{\text{th}} Y_j^{\text{th}} \sum_k \lambda_{ik} \lambda_{jk} \left(\frac{\Delta R_k}{R_k} \right)^2. \quad (19)$$

The uncertainties in the experimental data (1), (2), (4) should be added to the diagonal matrix elements of the error matrix (19)

$$\sigma_i^{\text{tot}2} = \sigma_{ii}^2 + \sigma_i^{\text{ex}2}. \quad (20)$$

For ${}^4\text{He}$ σ^{tot} differs from σ^{ex} insignificantly, while for D and especially for ${}^7\text{Li}$ σ_{ii} and σ_i^{ex} are comparable. If we neglect the correlations then the matrix w_{ij} is diagonal and equal to

$$w_{ii} = 1/\sigma_i^{\text{tot}2}.$$

In this case we can present the likelihood function (17) as a product of three individual functions $Lf(\eta, Q) = Lf_D(\eta, Q) Lf_{\text{He}}(\eta, Q) Lf_{\text{Li}}(\eta, Q)$ The equation

$$Y_i(\eta, Q) = Y_i^{\text{ex}} \quad (21)$$

defines three lines in the $\eta-Q$ plane where the individual likelihood functions are equal to one. The equations

$$[Y_i(\eta, Q) - Y_i^{\text{ex}}]^2 = \sigma_i^{\text{tot}2} \quad (22)$$

define 1σ ranges around these lines for each element. These ranges are shown in Fig. 2. The slope of the deuterium range is smaller than that of helium and lithium reflecting smaller sensitivity in Q and higher sensitivity in η .

In contrast, the helium range goes almost vertically reflecting high sensitivity of the helium fraction to Q and low sensitivity to η . This low sensitivity to η can be explained by a large helium binding energy. Only gamma's with the energy $E_\gamma > 20$ MeV can significantly change the number of helium nuclei. At any η the number of such γ quanta is small at the BBN temperature. We can, therefore, expect the low sensitivity of the helium mass fraction to η .

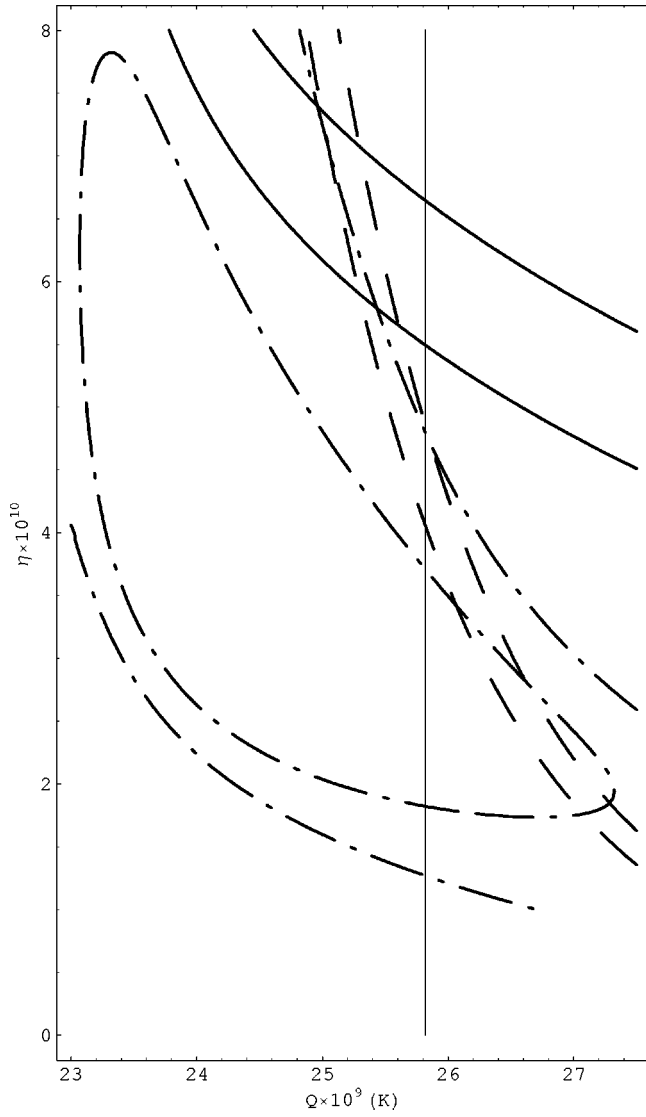


FIG. 2. 1σ ranges around the maxima of individual likelihood functions. The solid lines show 1σ ranges for D, the dashed lines are for ${}^4\text{He}$ [using Y_p from Eq. (5)], and the dot-dashed lines are for ${}^7\text{Li}$. For lithium, there are 2 solutions for η and Q , hence the shape of the error contours is more complicated.

The lithium range has two distinct branches corresponding to two different solutions of Eq. (21) for η at given Q . All three ranges intersect near $\eta=6.5$ and $Q=25$. One can expect that the general likelihood function (17) will have a maximum in this region. Indeed, we found the maximum of $Lf(\eta, Q)$ at the point $\eta_m = (6.51 + 0.77 - 0.66) \times 10^{-10}$ and $Q_m = (25.26 \pm 0.20) \times 10^9$ K. Figure 3 shows 1σ elliptic boundary near the maximum. The long axis of the ellipsis is almost vertical. Therefore, the correlation between $\Delta\eta$ and ΔQ is not significant. Comparing Figs. 3 and 2 one can conclude that the error ΔQ is determined mostly by ${}^4\text{He}$ mass fraction data. It is interesting to note that η_m is compatible with the one found from recent CMB anisotropy measurements [43]. The dark shadow region shows the 1σ range for η fitted from BBN only at present value of $Q = 25.82$ K.

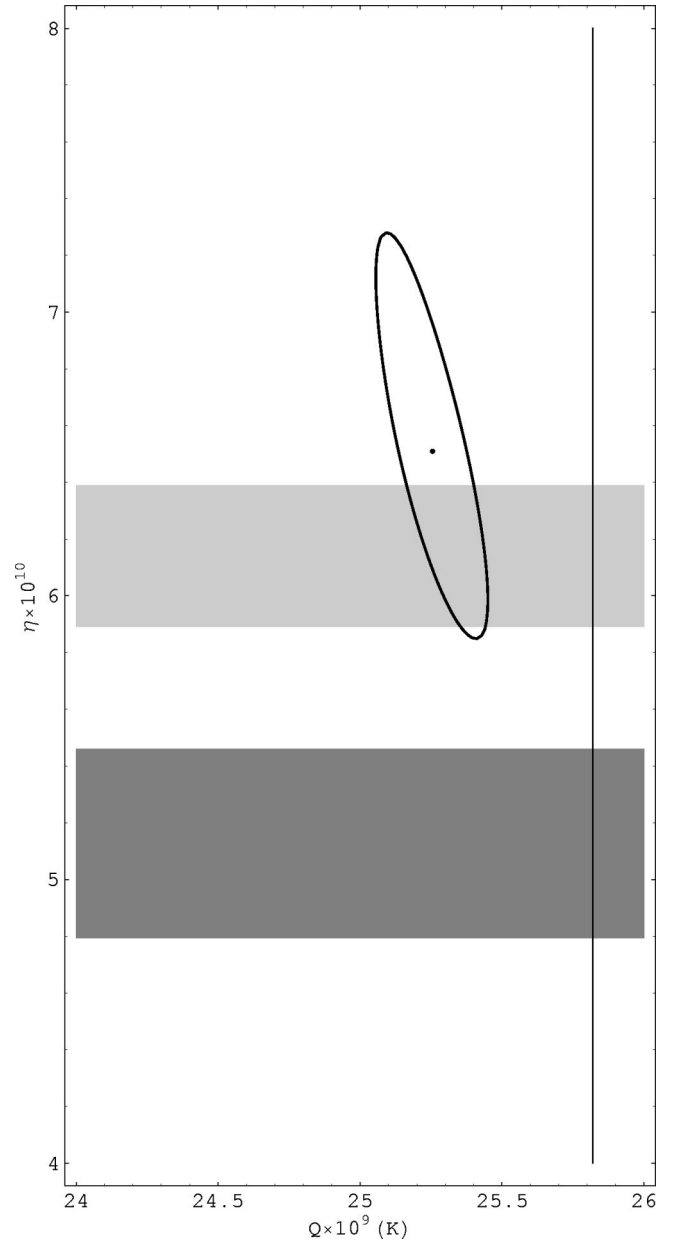


FIG. 3. 1σ range about the maximum of $Lf(\eta, Q)$ [again using Y_p from Eq. (5)]. The lighter shaded region shows CMB-WMAP data for η . The darker shaded region is the 1σ -range for η from BBN calculations using the present-day value of the deuteron binding energy, $Q = 25.82$. A lower value of Y_p will produce a larger deviation between the η_{WMAP} and η_{BBN} .

B. Constraint from CMB anisotropy measurements

The value of η found from CMB anisotropy measurements

$$\eta_0 = (6.14 \pm 0.25) \times 10^{-10}$$

has a rather high accuracy. It is natural to use the constraint from this measurement in our study of the deuteron binding energy effects. To do this we construct another likelihood function which is a function of Q only:

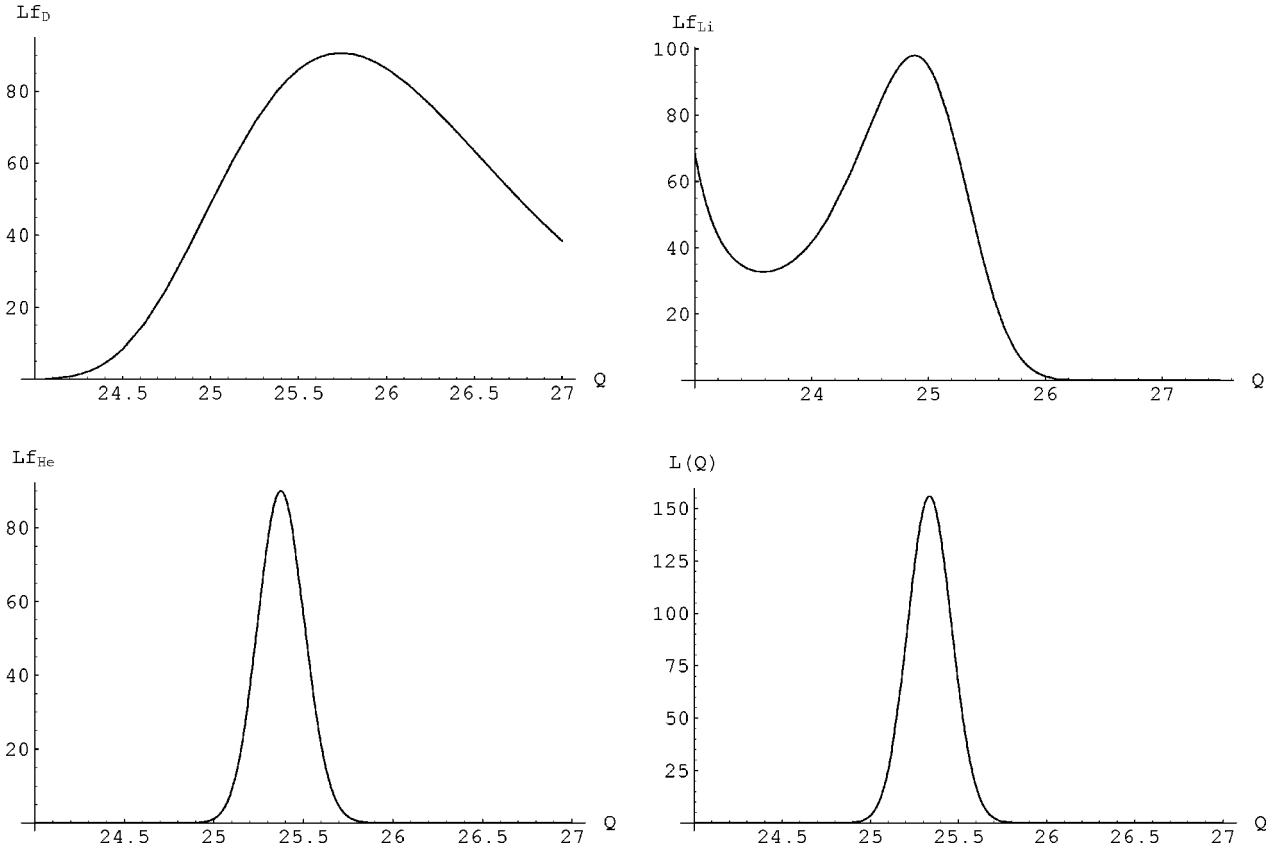


FIG. 4. Individual likelihood functions (23) for the light elements. From top to bottom: D, ${}^4\text{He}$ [Eq. (5)], Li, and the combined dataset.

$$L(Q) = \int_{-\infty}^{\infty} \exp\left(-\frac{(\eta - \eta_0)^2}{2\sigma_\eta^2}\right) Lf(\eta, Q) d\eta. \quad (23)$$

If we neglect nondiagonal elements in w_{ij} we can construct the individual likelihood functions for D, ${}^4\text{He}$, and ${}^7\text{Li}$. They are constructed in the same way as Eq. (23) using instead of general function $Lf(\eta, Q)$ the individual ones $Lf_D(\eta, Q)$, $Lf_{\text{He}}(\eta, Q)$, $Lf_{\text{Li}}(\eta, Q)$. These functions are plotted in Fig. 4 together with the general likelihood function (23).

From the deuterium likelihood function we found the position of the maximum and 1σ deviations

$$Q_D = (25.74 + 0.92 - 0.68) \times 10^9. \quad (24)$$

The shape near the maximum is apparently non-symmetric. The position of the maximum is fully compatible with the present value of $Q = 25.82 \times 10^9$ K. The helium likelihood function is much narrower (see the second panel from the top). It gives for the maximum and for the 1σ the values

$$Q_{\text{He}} = (25.37 \pm 0.13) \times 10^9. \quad (25)$$

This value lies below the present value of the binding energy. Finally, the lithium likelihood function has the maximum at

$$Q_{\text{Li}} = (24.88 + 0.43 - 0.59) \times 10^9. \quad (26)$$

The position of this maximum is compatible with the helium result.

The general likelihood function (23) is plotted in the lower panel in Fig. 4. The position of its maximum differs only slightly from the position given by the helium likelihood function:

$$Q_{\text{BBN}} = (25.34 \pm 0.12) \times 10^9. \quad (27)$$

It is interesting to compare the light element abundances for two values of the deuterium binding energies. In Fig. 5 we plotted the traditional curves for the light element abundances as a function of η for two values of Q . The dotted lines in the figures correspond to a present value of $Q_{\text{present}} = 25.82 \times 10^9$ K, while the solid curves correspond to a new value $Q_{\text{BBN}} = 25.34 \times 10^9$ K. Clearly, the new value Q_{BBN} moves the curves closer to the data.

The result which we obtained may be presented as

$$\delta Q/Q = -0.019 \pm 0.005, \quad (28)$$

where $\delta Q = Q_{\text{BBN}} - Q_{\text{present}}$. If we do not fix η and try to fit it simultaneously with Q we obtain

$$\delta Q/Q = -0.022 \pm 0.008, \quad \eta = (6.51 + 0.77 - 0.66) \times 10^{-10}. \quad (29)$$

The obtained η is fully compatible with the one measured by WMAP.

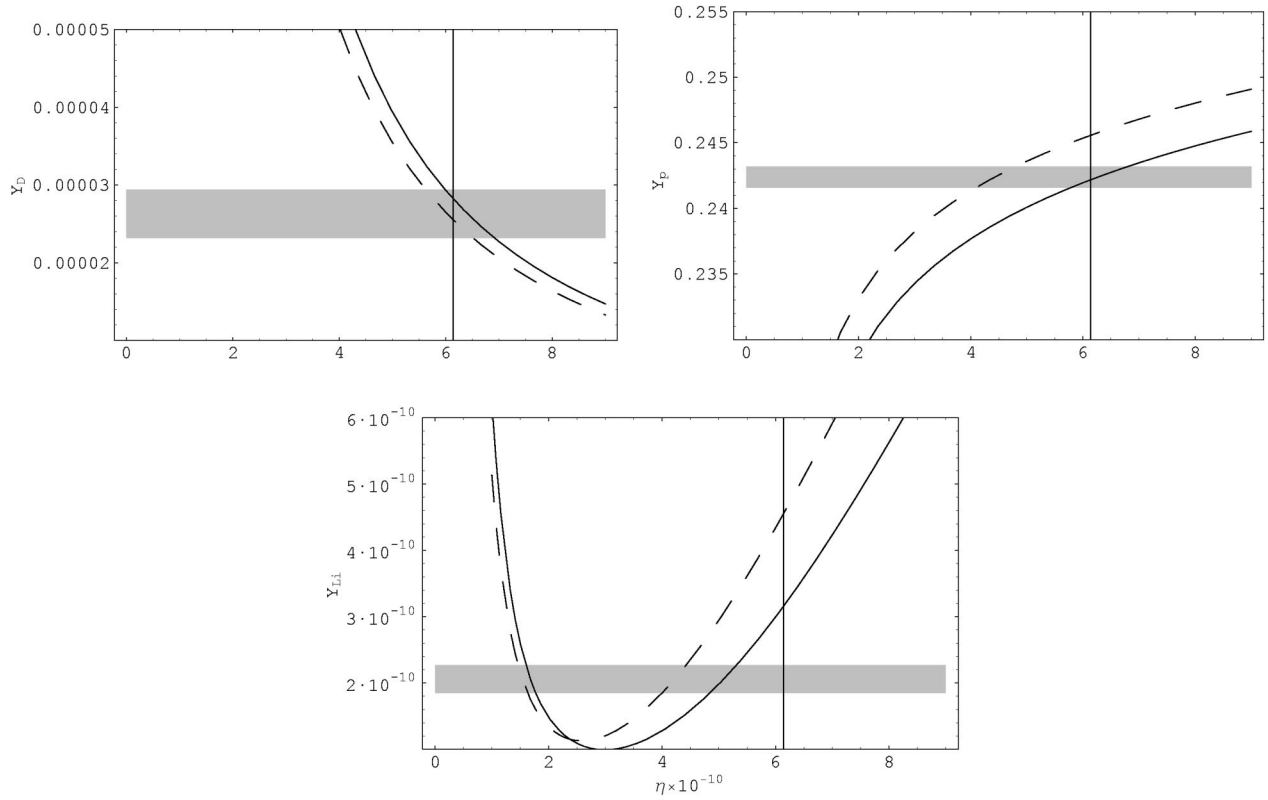


FIG. 5. The predicted light element abundance yields as a function of η , for two values of the deuteron binding energy Q . The dotted curve corresponds to the present value of $Q_{\text{present}} = 25.82 \times 10^9$ K. The solid curve corresponds to the new value of $Q = Q_{\text{BBN}} = 25.34 \times 10^9$ K. The vertical line corresponds to $\eta = 6.14$ (WMAP value). The shaded regions is the 1σ ranges for the observed light element abundances, where Y_p is from Eq. (5).

These values of $\delta Q/Q$ and η were obtained for high value of the helium mass fraction Y_p . If we use as an input the low value of Y_p from Eq. (6) we obtain

$$\delta Q/Q = -0.048 \pm 0.004. \quad (30)$$

If we fit both $\delta Q/Q$ and η we obtain

$$\delta Q/Q = -0.059 \pm 0.007, \quad \eta = (7.55 + 0.91 - 0.75) \times 10^{-10}. \quad (31)$$

Finally if we use the value of Y_p for ^4He obtained using the whole sample of 14 points, with increased error bars, from Eq. (1), we obtain

$$\delta Q/Q = -0.033 \pm 0.006, \quad (32)$$

and for $\delta Q/Q$ and η

$$\delta Q/Q = -0.042 \pm 0.009, \quad \eta = (7.00 + 0.85 - 0.72) \times 10^{-10}. \quad (33)$$

The results given in Eqs. (28) and (30) therefore represent an estimate of the plausible range in $\delta Q/Q$. Despite the clear systematic uncertainties in the ^4He data, and accepting the WMAP value of η as being correct, $\delta Q/Q$ appears to deviate from zero by 4σ [Eq. (28)] or greater [Eqs. (30), (32)].

The deuteron binding energy depends on the strong scale and quark masses. It is convenient to assume that Λ_{QCD} is

constant, and the quark mass is variable. This only means that we measure all energies in units of Λ_{QCD} (and cross sections in units $\Lambda_{\text{QCD}}^{-2}$). In Ref. [5] we concluded that the deuteron binding energy is very sensitive to variation of the strange quark mass m_s [45]:

$$\frac{\delta(Q/\Lambda_{\text{QCD}})}{(Q/\Lambda_{\text{QCD}})} = -17 \frac{\delta(m_s/\Lambda_{\text{QCD}})}{(m_s/\Lambda_{\text{QCD}})}. \quad (34)$$

Combining Eqs. (32) and (34) we obtain

$$\frac{\delta(m_s/\Lambda_{\text{QCD}})}{(m_s/\Lambda_{\text{QCD}})} = (1.1 \pm 0.3) \times 10^{-3}. \quad (35)$$

This equation may contain an additional factor (close to one) reflecting unknown theoretical uncertainty in Eq. (34). Note that we obtain here variation at the level 10^{-3} while the limits on variation of α [2,46] and $\Lambda_{\text{QCD}}/M_{\text{plank}}$ [2,3] are an order of magnitude weaker. This may serve as a justification of our approach.

IV. CONCLUSION

Allowing the deuteron binding energy Q to vary in BBN appears to provide a better fit to the observational light element abundance data. Varying Q simultaneously does two things; it resolves the internal inconsistency between ^4He

and the other light elements, and it also results in excellent independent agreement with the baryon to photon ratio determined from WMAP (Fig. 5). However, the magnitude of the variation is sensitive primarily to the observed ${}^4\text{He}$ abundance, which has the smallest relative statistical error. A systematic error in the abundance of ${}^4\text{He}$ could imitate the effect of the deuteron binding energy variation, although one needs a systematic error which is very much greater than has been claimed in the most recent observational work.

We note that Izotov and Thuan [27], the most recent estimate for Y_p in our sample, argue that systematics are at most 0.6% for that survey. On the other hand, the possibility has also been explored that the creation of ${}^4\text{He}$ in population III stars might mean that the true primordial ${}^4\text{He}$ abundance is lower even than that seen in the most metal-poor objects [47]. If so, the significance of the deviation of $\delta Q/Q$ from zero we report in this paper would be even larger. These results hopefully provide an extremely strong motivation to obtain substantially better measurements of all the light elements, and to explore even more intensively, the possible sources of systematic errors.

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APPENDIX

Let $U_0(r)$ be a critical depth potential for which the binding energy $Q=0$, and $U_t(r)$ a potential for a proton neutron system in a triplet state producing a deuteron with small binding energy $Q=2.22$ MeV. If we add to the deuteron Hamiltonian a perturbation

$$\delta U_\lambda(r) = \lambda[U_0(r) - U_t(r)],$$

then, variation of λ from 0 to 1 will move the binding energy Q from 2.22 MeV to 0. From a virial theorem for a quantum system we have

$$\frac{dE}{d\lambda} = \int_0^\infty [U_0(r) - U_t(r)]\chi^2(r)dr, \quad (\text{A1})$$

where $\chi(r)$ is the radial s-wave function. For simplicity we neglect the d-wave contribution. For $Q \rightarrow 0$ the main contribution into normalization integral for $\chi(r)$ comes from the

region outside of the nuclear forces radius R . The normalization integral can be presented as sum of contributions from inner and outer regions

$$\int_0^R \chi^2(r)dr + b^2 \int_R^\infty e^{-2\gamma r} dr = 1, \quad (\text{A2})$$

where $\gamma = \sqrt{m_p|E|/\hbar^2}$. At $|E| \rightarrow 0$ the second integral dominates giving $b^2 = 2/\gamma$. Separating the E dependence of the normalization factor in $\chi(r)$ we can rewrite Eq. (A1) as

$$\frac{dE}{\sqrt{|E|}} = d\lambda 2 \sqrt{\frac{m_p}{\hbar^2}} \int_0^\infty [U_0(r) - U_t(r)]\tilde{\chi}^2(r)dr, \quad (\text{A3})$$

where $\tilde{\chi}(r)$ is practically independent on E inside the potential well (where $E \ll U$) and $\tilde{\chi}(r) = e^{-2\gamma r} \rightarrow 1$ at $r > R$ when $|E| \rightarrow 0$. Integrating the left hand side of Eq. (A3) over E from $-Q$ to 0 and the right hand side of Eq. (A3) over λ from 0 to 1 we obtain

$$Q = \frac{m_p}{\hbar^2} \left(\int_0^\infty [U_0(r) - U_t(r)]\tilde{\chi}^2(r)dr \right)^2. \quad (\text{A4})$$

Equation (A4) shows that the position of a shallow bound level depends quadratically on the difference between the actual depth of the potential and the critical one. For a square well $Q = \pi^2(U - U_0)^2/16U_0$, $U_0 = \pi^2\hbar^2/4m_p R^2$.

In fact, the Eq. (A4) is valid not only for the energy of a bound level but for the energy of a virtual level as well. The integration in Eq. (A4) is over the region $r < R$ where the function $\tilde{\chi}^2(r)$ is insensitive to the energy E , and the quadratic dependence on $U - U_0$ guarantees the validity of Eq. (A4) for both $U < U_0$ and $U > U_0$. Thus, for the energy of the virtual level we have

$$\epsilon_v = \frac{m_p}{\hbar^2} \left(\int_0^\infty [U_0(r) - U_s(r)]\tilde{\chi}^2(r)dr \right)^2, \quad (\text{A5})$$

where $U_s(r)$ is the potential for a singlet states. We have both $Q \ll U$ and $\epsilon_v \ll U$. This means that the difference between the triplet and singlet potentials is not large. Assuming that the changes in the triplet and singlet potentials are the same we obtain for the changes in Q and ϵ_v the relation

$$\frac{\delta\epsilon_v}{\sqrt{\epsilon_v}} = -\frac{\delta Q}{\sqrt{Q}}. \quad (\text{A6})$$

This equation also holds for the effect produced by variation of the proton mass (the dominating effect comes from variation of U_0).

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