

P-state positronium for precision physics: An ultrafine splitting at α^6

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An “ultrafine” splitting in positronium between the $L = 1$ spin-singlet state and the spin average of the spin-triplet states is shown to arise only at order α^6 . The QED prediction for $n = 2$ states is $\Delta_{2,P} = (683/172\,800)m\alpha^6 = 73.7(2.6)$ kHz. This represents the smallest leading-order QED splitting known. Current experimental efforts could observe this splitting, and its observation can constrain new ultralight interactions, such as axions or Z' .

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I. INTRODUCTION

Spectroscopy of positronium represents a precision test of quantum electrodynamics (QED). The low mass of the electron compared to other mass scales renders the contributions from strong, electroweak, and heavier lepton loops negligible within the accuracy of current experiments. QED predictions for the positronium energy levels exist completely at order α^6 [1–5]. Research is active to compute the incomplete α^7 correction [6–21]. These predictions are in agreement with existing measurements for a number of level transitions: $2^3S_1-1^3S_1$ ($1S$ - $2S$ interval) [22], $1^3S_1-1^1S_0$ (ground-state hyperfine splitting) [23,24], and $2^3S_1 \rightarrow 2^{2S+1}P_J$ (fine structure and Lamb shift) [25,26].

Energy levels are derived as a double power series in $\alpha^g \ln^h(\alpha)$ starting at α^2 . The coefficients $C_{gh}(n, L, J, S)$ can be zero [i.e., the α^3 and $\alpha^4 \ln^h(\alpha)$ terms]. Cancellations between $C_{gh}(n, L, J, S)$ can further occur in level transitions. The $1S$ - $2S$ interval depends only on spin-independent coefficients and begins at α^2 . The fine-structure and ground-state hyperfine splittings (hfs’s) are sensitive only to spin-dependent terms and are nonzero at α^4 . The Lamb shift (defined as the spin-independent $2S$ - $2P$ splitting) starts only at $\alpha^5 \ln(\alpha)$.

This work focuses on the hyperfine splitting between the spin-singlet state n^1P_1 and the spin average of the spin-triplet n^3P_J states:

$$\Delta_{n,P} \equiv M(n^1P_1) - \frac{1}{9}[M(n^3P_0) + 3M(n^3P_1) + 5M(n^3P_2)], \quad (1)$$

where $M(n^{2s+1}L_J)$ is the energy level of the state of given quantum numbers. Naively, one might expect this splitting to be on the order of gigahertz like the Lamb shift or hfs. Using the existing fine-structure measurements (compiled in Table I), one instead finds $\Delta_{2,P} = 4.31(6.50)$ MHz [26–28]. This result is consistent with zero within the experimental uncertainty at the parts per thousand level. Analogous experimental splittings in heavy quarkonium are consistent with zero to even higher precision, and Eq. (1) has been suggested as a method for studying the exotic spectrum of quarkonium [29].

In this paper, we investigate the physical origin underpinning this precise cancellation in positronium. We show that this particular hyperfine splitting arises only at α^6 and is further

suppressed parametrically, justifying calling it an *ultrafine* splitting. Finally, we briefly discuss how $\Delta_{n,P}$ constrains new ultralight interactions.

II. ULTRAFINE SPLITTING

To understand the smallness of $\Delta_{n,P}$, it is useful to study the matrix elements of operators contributing to a $f\bar{f}$ bound state following the same discussion in the context of heavy quarkonium [29]. The set of spin-independent operators depending on powers of squared momenta is infinite. In stark contrast, only a finite set of linearly independent spin-dependent operators exists given the fermion spins $\mathbf{S}_f, \mathbf{S}_{\bar{f}}$ and the orbital angular momentum \mathbf{L} . Wigner-Eckart theorem restricts nonzero matrix elements of spin operators in states with total spin S to those which transform under an irreducible representation $k < 2S$. For states built from only two fermions, $S = 0, 1$, and therefore the largest irreducible representation is $k = 2$.

A particularly useful set of linearly independent matrix elements up to quadratic order is

$$\mathbf{S}_f \cdot \mathbf{S}_{\bar{f}} \quad (\text{hyperfine}), \quad (2)$$

$$\mathbf{S} \cdot \mathbf{L} \quad (\text{spin orbit}), \quad (3)$$

$$\vec{T} \equiv (\mathbf{S}_f \cdot \hat{r})(\mathbf{S}_{\bar{f}} \cdot \hat{r}) - \frac{1}{3}\mathbf{S}_f \cdot \mathbf{S}_{\bar{f}} \quad (\text{tensor}), \quad (4)$$

where $\mathbf{S} \equiv \mathbf{S}_f + \mathbf{S}_{\bar{f}}$. In a given spin multiplet, all other spin operators can be constructed from this set.

Using Wigner-Eckart theorem, the spin-orbit and tensor matrix elements must vanish for $S = 0$ states because they transform as $S = 1$ and $S = 2$, respectively. For $L = 0$ states, these matrix elements are also zero, so any $n^3S_1-n^1S_0$ splitting depends only on $\mathbf{S}_f \cdot \mathbf{S}_{\bar{f}}$. For states of $S = 1$ and $L > 0$, the spin-orbit and tensor matrix elements do not vanish but are given via the total angular momentum $J = L + 1, L$, and $L - 1$:

$$\langle \mathbf{S} \cdot \mathbf{L} \rangle = \frac{1}{2}[J(J+1) - L(L+1) - S(S+1)] \quad (5)$$

and

$$\langle \vec{T} \rangle = \begin{cases} -\frac{L+1}{6(2L-1)}, & J = L - 1, \\ +\frac{1}{6}, & J = L, \\ -\frac{L}{6(2L+3)}, & J = L + 1. \end{cases} \quad (6)$$

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TABLE I. Experimental and theoretical values for the $n = 2$ fine-structure transitions. The theoretical error is obtained from estimating the size of higher-order corrections and is explained in the text with more detail.

Transition	ΔE_{exp} (MHz)	ΔE_{theory} (MHz)	δE (MHz)
$2^3S_1 \rightarrow 2^1P_1$	11180(5)(4) [25]	11185.37(8) [1]	-5.37(500)(400)(8)
$2^3S_1 \rightarrow 2^3P_0$	18499.65(120)(400) [26]	18498.25(8) [1]	1.40(120)(400)(8)
$2^3S_1 \rightarrow 2^3P_1$	13012.42(67)(154) [26]	13012.41(8) [1]	-0.01(67)(154)(8)
$2^3S_1 \rightarrow 2^3P_2$	8624.38(54)(140) [26]	8626.71(8) [1]	2.30(54)(140)(8)

In this multiplet, $(\mathbf{S} \cdot \mathbf{L})^2$ occurs as an order α^6 correction, and therefore the following relation is useful [30]:

$$\langle (\mathbf{S} \cdot \mathbf{L})^2 \rangle = \frac{L(L+1)}{3} \langle \mathbf{S}^2 \rangle - \langle \mathbf{S} \cdot \mathbf{L} \rangle - (2L-1)(2L-3) \langle \vec{T} \rangle, \quad (7)$$

where \mathbf{S}^2 has been used for compactness but is directly related to $\mathbf{S}_f \cdot \mathbf{S}_{\bar{f}}$.

Due to these additional nonzero matrix elements in $L > 0$ states, individual $2^3S_1 \rightarrow 2^{2S+1}P_J$ transitions have a more complicated spin dependence. The spin-orbit and tensor elements can be made to vanish by construction in a sum weighted by the $2J+1$ degenerate spin states. This sum receives contributions only from $\mathbf{S}_f \cdot \mathbf{S}_{\bar{f}}$ and is given by

$$\begin{aligned} \Delta_{n,L} \equiv & M(n^1L_{J=L}) - \frac{2L-1}{3(2L+1)} M(n^3L_{J=L-1}) \\ & - \frac{2L+1}{3(2L+1)} M(n^3L_{J=L}) - \frac{2L+3}{3(2L+1)} M(n^3L_{J=L+1}), \end{aligned} \quad (8)$$

from which we have derived Eq. (1) for the $L = 1$ case. $\Delta_{n,L}$ can be thought of as analogous to the hfs for $L > 0$ states as both are purely hyperfine operator dependent. One might guess that $\Delta_{n,L}$ should be an order α^4 observable like the hfs. Instead, due to the dynamics underlying the nonrelativistic hyperfine operator, $\Delta_{n,L}$ arises only at higher order in positronium.

The nonrelativistic positronium potential $V(\mathbf{r})$ to order α^5 has long been known (a modern derivation using nonrelativistic QED can be found in Ref. [31]). The hyperfine operator arises at α^4 , and the only term up to α^5 is proportional to $\delta^{(3)}(\mathbf{r})$ [32], implying it arises as contact interaction from the reduction of the QED interaction. A natural way to see this starts from the Dirac equation with the Breit interaction [33]. In this Hamiltonian, the spin-spin coupling comes from the Laplacian operator acting on the Coulomb potential $1/r$, i.e., the Fourier transform of the massless propagator $1/q^2$. This term is a generic feature of massless gauge theories like quantum chromodynamics as well and therefore a source of hyperfine splitting in quarkonium.

The nonrelativistic wave functions near the origin scale as r^L ; thus only $L = 0$ states receive contributions from $\delta^{(3)}(\mathbf{r})$ terms. Therefore $\Delta_{n,L}$ must come from at least α^6 corrections with more nontrivial r dependence. The vanishing of $\Delta_{n,L}$ at $O(\alpha^5)$ is a unique characteristic of positronium. Heavier leptonic systems like true muonium ($\mu^+\mu^-$) have typical momentum $\alpha m_\mu \sim m_e$ and therefore receive corrections at α^5 due to electron loop corrections to the potential [34,35].

For quarkonium states, both lighter flavors and the gluon self-coupling introduce corrections at α_s^5 [29].

For the $L = 1$ states, the α^6 contributions were computed in Refs. [3,4] (with small misprints corrected in Ref. [1]). The general $L > 0$ state energy levels up to order α^6 were computed in Ref. [5]. From these, we find that the nonzero contributions to $\Delta_{n,P}$ come from

$$\begin{aligned} \delta E = & \frac{m\alpha^6}{n^3} \left[\frac{1}{2000} \left(46 - \frac{43}{n^2} \right) (\mathbf{S} \cdot \mathbf{L})^2 - \frac{1}{2000} \left(14 - \frac{17}{n^2} \right) \mathbf{S}^2 \right. \\ & \left. - \frac{1}{7680} \left(227 + \frac{90}{n} - \frac{108}{n^2} \right) \kappa^2 \right], \end{aligned} \quad (9)$$

where $\kappa = \frac{1}{5}[-2(\mathbf{L} \cdot \mathbf{S})^2 + 4\mathbf{L} \cdot \mathbf{S} + \frac{4}{3}\mathbf{S}^2]$ and terms known to vanish in $\Delta_{n,P}$ have been neglected. The remaining terms are all necessarily proportional to $\mathbf{S}_f \cdot \mathbf{S}_{\bar{f}}$. This expression is the sum of a number of physically distinct processes: second-order relativistic corrections to the Coulomb potential, first-order relativistic corrections to single magnetic photon exchange, and iterating the Breit interaction [3,4]. These contributions are nonzero because they arise from higher moments of the wave function $\langle \mathbf{r}^{-k} \rangle$.

Using Eq. (7), we can derive a general relation for reducing general spin-operator contributions into $\mathbf{S}_f \cdot \mathbf{S}_{\bar{f}}$ ones:

$$\begin{aligned} \Delta_{n,L} \langle (\mathbf{S}^2)^m (\mathbf{L} \cdot \mathbf{S})^n \rangle \\ = \frac{2^m}{9} [(-1-L)^n + 3(-1)^n + 5L^n] \Delta_{n,L} \langle \mathbf{S}_f \cdot \mathbf{S}_{\bar{f}} \rangle, \end{aligned} \quad (10)$$

where $\Delta_{n,L}(\mathcal{O})$ defines the spin-averaged matrix element of the operator \mathcal{O} in the spin-triplet with quantum numbers n and L . An expression for $\Delta_{n,L>1}$ could be derived by using the results of [5], but we see that $\Delta_{n,L>1} \propto \alpha^6$ as well. These splittings will be even more parametrically suppressed due to the larger n required, but given that no experimental measurements have been undertaken to measure $L > 1$ state energies, a nonzero measurement of $\Delta_{n,L>1}$ is not possible in the foreseeable future. For P states, we then find that

$$\Delta_{n,P} = \frac{m\alpha^6}{60} \left(\frac{137}{90n^3} + \frac{1}{n^4} - \frac{1}{2n^5} \right) \Delta_{n,P} \langle \mathbf{S}_f \cdot \mathbf{S}_{\bar{f}} \rangle. \quad (11)$$

Using $\Delta_{n,P} \langle \mathbf{S}_f \cdot \mathbf{S}_{\bar{f}} \rangle = -1$ and setting $n = 2$, we find

$$\Delta_{2,P} = \frac{683m\alpha^6}{172800} = 73.7(2.6) \text{ kHz}, \quad (12)$$

which is about 90 times smaller than the existing experimental accuracy. The smallness of $\Delta_{2,P}$ compared to other splittings is dramatic. The hfs is 203 GHz, while the typical Lamb-shift-like splittings in Table I are 10 GHz. The theoretical error was

estimated by considering the α expansion,

$$\Delta_{n,P} = m\alpha^6 \left(C_0 + C_{11}\alpha \ln \alpha + C_{10} \frac{\alpha}{\pi} + \dots \right), \quad (13)$$

where the coefficients C_{gh} start at α^6 .

There are a number of ways to estimate the error; one is to take the leading unknown contribution with a coefficient of 1. This would be 670 kHz for the missing $\alpha^7 \ln \alpha$ to $\Delta_{2,P}$. A further assumption could be to take n^{-3} scaling to yield an estimate of 80 kHz. Another common estimate is to take some fraction of the highest-order contribution that is known. This was done previously for the values in Table I, where the error is given by half the $\alpha^7 \ln^2(\alpha)$ contribution [1]. This would suggest an estimate of 40 kHz for $\Delta_{2,P}$. The knowledge that the coefficients of $\Delta_{2,P}$ are parametrically suppressed both because $|\psi_{nP}(0)| \sim 0$ and because of the cancellations by construction, these estimates are unreasonably conservative, and we propose another. We take $C_0 \sim C_{11} \sim C_{10}$ to approximate the parametric suppression of the coefficients, in which case the $O(\alpha \ln \alpha)$ term dominates the theoretical error in Eq. (12). Comparing this to Table I, we see removing the dependence upon 2^3S_1 where there is no parametric suppression reduces the theoretical error by a factor of 30.

Having shown that existing measurements are too imprecise to resolve this splitting, one can ask about the future prospects. Upcoming experiments at University College London plan to remeasure the $n = 2$ fine structure with a reduced uncertainty, yielding $\Delta_{2,P} \sim 100$ kHz [36–38]. A measurement of $\Delta_{2,P}$ would require only a factor of 2 improvement. A precision of ~ 10 kHz would translate into a 5σ detection.

The spin weighting of the ultrafine splitting means the uncertainty on $2^3S_1 \rightarrow 2^1P_1$ contributes 86% of the error, assuming the transitions are measured with equal accuracy. The situation is worse because $2^3S_1 \rightarrow 2^1P_1$ is a forbidden single-photon transition. Previous measurements of this transition were performed in magnetic fields where mixing

with the other $2P$ states occurs. With this mixing, the single-photon transition is allowed, and the multiple measurements at different magnetic fields are extrapolated to the zero-field limit via the known theoretical dependence [39]. This method inherently has larger uncertainties compared to the allowed $2^3S_1 \rightarrow 2^3P_J$ states.

Beyond measuring QED precisely, the ultrafine splitting can constrain new ultralight spin-dependent interactions like the exchange of an axion or Z' . The nonrelativistic reduction of these interactions is a Yukawa potential $\alpha' \mathbf{S}_f \cdot \mathbf{S}_{\bar{f}} \frac{e^{-m_\phi r}}{r}$, where α' is the new coupling and m_ϕ is the mass of the new particle. From this, the correction to $\Delta_{2,P}$ is

$$\delta\Delta_{2,P} = \frac{m\alpha\alpha'}{8\left(1 + 2\frac{m_\phi}{m\alpha}\right)^4}. \quad (14)$$

Comparing this expression to Eq. (12), we see that a measurement of $\Delta_{2,P}$ constrains $\alpha' \sim \alpha^5 \sim 10^{-11}$ in the $m_\phi = 0$ limit. This would be competitive with other model-independent constraints on ultralight particles in atomic systems [40–45].

III. CONCLUSIONS

In this work, we have pointed out a previously overlooked ultrafine splitting in positronium. At $O(m\alpha^6)$, the $\Delta_{2,P} = 0.0737(12)$ MHz splitting represents the smallest leading-order splitting from QED known. While existing measurements are too imprecise, it should be resolved in the near term. Further theoretical work needs to be done to compute the $O(m\alpha^7 \ln \alpha)$ contribution. Additionally, this very small effect presents a new opportunity to constrain ultralight particles with coupling to electrons.

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