## Status of QED prediction of the electron g-2

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• current experimental value of the anomaly

• current status of the theoretical calculation

• recent results (mostly by S. Laporta)

if  $\vec{\mu}$  is the magnetic moment of the electron and  $\vec{s}$  its spin

$$\vec{\mu} = g_e \frac{e\hbar}{2mc} \vec{s}$$

where  $g_e$  is the gyromagnetic ratio; the Dirac equation gives  $g_e = 2$  for a free electron – but  $g_e \neq 2$  (Breit) for a *bound* electron ( $g_e$ -bound), so in general

 $g_e = 2(1 + a_e)$ 

where  $a_e = \frac{1}{2}(g-2)$  is the anomaly. in QED,  $g_e \neq 2$ , or  $a_e \neq 0$  even for the *free* electron. Experimental values of the *free* electron  $a_e$  (current main problem: cavity shift)  $10^{12} a_e(\exp) = 1\,159\,652\,188.4\,(4.3)[3.7 \text{ ppb}]$  Dehmelt 1987,UW  $= 1\,159\,652\,180.85(.76)[0.66 \text{ ppb}]$  Gabrielse 2006, Harvard  $= 1\,159\,652\,180.73(.28)[0.24 \text{ ppb}]$  Gabrielse 2008, Harvard PRL 100, 120801 (2008)

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## G

## New Measurement of the Electron Magnetic Moment and the Fine Structure Constant

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A measurement using a one-electron quantum cyclotron gives the electron magnetic moment in Bohr magnetons,  $g/2 = 1.001\,159\,652\,180\,73\,(28)$  [0.28 ppt], with an uncertainty 2.7 and 15 times smaller than for previous measurements in 2006 and 1987. The electron is used as a magnetometer to allow line shape statistics to accumulate, and its spontaneous emission rate determines the correction for its interaction with a cylindrical trap cavity. The new measurement and QED theory determine the fine structure constant, with  $\alpha^{-1} = 137.035\,999\,084\,(51)$  [0.37 ppb], and an uncertainty 20 times smaller than for any independent determination of  $\alpha$ .

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The electron magnetic moment  $\mu$  is one of the few measurable properties of one of the simplest of elementary particles—revealing its interaction with the fluctuating QED vacuum, and probing for size or composite structure not yet detected. What can be accurately measured is g/2, the magnitude of  $\mu$  scaled by the Bohr magneton,  $\mu_B = e\hbar/(2m)$ . For an eigenstate of spin **S**,

$$\boldsymbol{\mu} = -\frac{g}{2} \mu_B \frac{\mathbf{S}}{\hbar/2},\tag{1}$$

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order QED calculation (underway [7]) than by the measurement uncertainty in g/2. The accuracy of the new gsets the stage for an improved CPT test with leptons. It also will allow an improved test of QED, and will be part of the discovery of low-mass dark-matter particles or the elimination of this possibility [8], when a better independent measurement of  $\alpha$  becomes available.

Figure 3 represents the lowest cyclotron and spin energy levels for an electron weakly confined in a vertical magnetic field  $B\hat{z}$  and an electrostatic quadrupole potential.

the theoretical expression for  $a_e$  is

$$a_{e} = a_{e}(QED)$$
  
+ $\Delta a_{e}(\mu) + \Delta a_{e}(\tau) + \Delta a_{e}(hadronic) + \Delta a_{e}(weak)$   
+ $\Delta a_{e}(NP)$ 

small terms:

$$\Delta a_e(\mu) = 2.71 \times 10^{-12}$$
$$\Delta a_e(\tau) = 0.01 \times 10^{-12}$$
$$\Delta a_e(\text{hadronic}) = 1.671(19) \times 10^{-12} \simeq \left(\frac{m_e}{m_\mu}\right)^2 \Delta a_\mu(\text{hadronic})$$
$$\Delta a_e(\text{weak}) = 0.030(01) \times 10^{-12}$$

The constraint on NP depends on the experimental error – but the dependence on unknown new particles is quadratic in the ratio of the masses (for large masses)

$$\Delta a_e(\mathrm{NP}) \simeq \left(\frac{m_e}{m_\mu}\right)^2 \Delta a_\mu(\mathrm{NP})$$

experimental error on  $a_e$ :  $\Delta a_e(\exp) = 0.28 \times 10^{-12}$ experimental values of the  $\mu$  anomaly

 $10^{12} a_{\mu}(\exp) = 1\ 165\ 920\ 800(600)[517\ ppb]$  E821 Brookhaven, 2004

$$\frac{\Delta a_{\mu}(\exp)}{\Delta a_{e}(\exp)} = 2100 \qquad \qquad \left(\frac{m_{\mu}}{m_{e}}\right)^{2} = 43000$$

the  $\mu$  anomaly is still more constraining on NP by a factor 20.

 $a_e(\texttt{QED})$  is the QED contribution with electron loops only and depends only on the fine structure constant  $\alpha$ 

$$a_e(QED) = C_1\left(\frac{\alpha}{\pi}\right) + C_2\left(\frac{\alpha}{\pi}\right)^2 + C_3\left(\frac{\alpha}{\pi}\right)^3 + C_4\left(\frac{\alpha}{\pi}\right)^4 + C_5\left(\frac{\alpha}{\pi}\right)^5 + \dots$$



 $= -0.328 478 965 579 \ldots$ , (Petermann, Sommerfield 1957) 7 diagrams

3 loops:

$$C_{3} = \frac{83}{72}\pi^{2}\zeta(3) - \frac{215}{24}\zeta(5) + \frac{100}{3}\left[\left(a_{4} + \frac{1}{24}\ln^{4}2\right) - \frac{1}{24}\pi^{2}\ln^{2}2\right]$$
  
$$- \frac{239}{2160}\pi^{4} + \frac{139}{18}\zeta(3) - \frac{298}{9}\pi^{2}\ln 2 + \frac{17101}{810}\pi^{2} + \frac{28259}{5184}$$
  
$$= 1.181\ 241\ 456\ \dots, \quad (S.Laporta, E.R.\ 1996) \quad 72\ diagrams$$

$$\zeta(p) = \sum_{n=0}^{\infty} \frac{1}{n^p}, \qquad a_4 = \sum_{n=0}^{\infty} \frac{1}{2^n n^4},$$
$$C_3 \left(\frac{\alpha}{\pi}\right)^3 = 14\ 804.20 \times 10^{-12} \ , \qquad \qquad \Delta a_e(\exp) = 0.28 \times 10^{-12}$$









891 diagrams

The 4-loop coefficient is known only numerically. Only a few diagrams are known in analytical form. Numerical values are obtained by using MonteCarlo integration of huge 11-dimensional integrands (... and fixing minor bugs underway)

$$C_4 = -1.5098(384)$$
 (Kinoshita, 1999)

$$C_4 = -1.7283(35)$$
 (Kinoshita, 2003) shift

 $C_4 = -1.9144(35)$  (Kinoshita, 2007)

shift of -0.24 due to the discovery of one error  $% \left( {{{\mathbf{r}}_{\mathbf{r}}}_{\mathbf{r}}} \right)$ 

shift of -0.22 due to the discovery of another error

$$C_4 \left(\frac{\alpha}{\pi}\right)^4 = (55.73 \pm 0.10) \times 10^{-12}$$
,  $\Delta a_e(\exp) = 0.28 \times 10^{-12}$ 

the numerical evaluation of the 5 loop contribution is underway (Kinoshita) guess:  $C_5 = 0.0 \pm 4$ 

$$C_5 \left(\frac{\alpha}{\pi}\right)^5 = (0 \pm 0.27) \times 10^{-12}$$
,  $\Delta a_e(\exp) = 0.28 \times 10^{-12}$ 

same order of magnitude as the current experimental error !

QED=QED(
$$\alpha$$
)  
QED +  $\alpha \rightarrow a_e$   
QED +  $a_e \rightarrow \alpha$   
QED +  $a_e + \alpha \rightarrow$  real check

best independent experimental values of  $\alpha$  from

$$\alpha^2 = \frac{2R_{\infty}}{c} \left(\frac{m_u}{m_e}\right) \left(\frac{M_X}{m_u}\right) \left(\frac{h}{M_X}\right)$$

where  $m_u$  is the unit of mass,  $M_X$  the mass of the atom X and  $(h/M_X)$  is measured by recoil spectroscopy

best results

 $\alpha^{-1}(Cs) = 137.036\ 000\ 00(110)$  (8.0 ppb) NIST, 2006  $\alpha^{-1}(Rb) = 137.035\ 998\ 78\ (91)$  (6.7 ppb) ENS, 2006

much larger than the (0.24 ppb) error in  $a_e(\exp)$ ; (relatively) weak check of QED by combining available theory and the Harvard 2008 value of  $a_e$ 

$$\alpha^{-1}(a_e) = 137.035\ 999\ 084\ (33)(39)$$
  
= 137.035\ 999\ 084\ (51) [0.37 ppb]

 $a_e$  and  $\alpha$ 



FIG. 1. Most accurate measurements of the electron g/2 (a), and most accurate determinations of  $\alpha$  (b).

We (mostly S. Laporta) are trying to improve  $C_4$ 

- to provide a check ;
- to reduce the error.

The by now standard approach (implemented by S.L. in his computer programs) is:

- Extraction of the contribution to g-2 from each diagram as a sum of scalar loop integrals.
- Reduction of each contribution to the combination of a (small) number of Master Integrals.
- high-precision numerical calculation of the Master Integrals.
- Cross-checks of procedures, intermediate and final results

the first step is almost trivial (with an algebraic program like, say,  $\tt SCHOONSCHIP$  or  $\tt FORM)$ 

The scalar loop integrals are reduced to Master Integrals by exploiting the integration by parts identities (Chetyrkin, Tkachov 1981) in the *D*-continuous regularization scheme.

Not surprisingly, the Laporta algorithm is used for solving the identities:

- an essential point of the algorithm is the *ordering* of the scalar integrals;
- a *sufficiently* large system of identities is built, containing all the relevant scalar integrals and more identities than integrals;
- as a rule, a large linear system of identities is obtained, up to  $10^6 \dots 10^7$  identities for graph.

• the system is then solved, one equation at the time, with the Gauss substitution rule, by expressing "more difficult" integrals (within the *ordering rule*) in terms of "simpler" ones;

As a result all the integrals are reduced to a combination of irreducible master integrals  $M_j$ 

$$a_e(\texttt{diagram}) = \sum_{j=1}^N \frac{p_j(D)}{q_j(D)} M_j(D)$$

where  $r_j(D), s_j(D)$  are polynomials in the dimension D with integer coefficients.

• at 3-loop, N = 18, *i.e.* the final result from all the graphs is expressed in terms of 18 *Master Integrals* only;

 $\bullet$  at 4-loop,  $N\sim 300$  is expected.

some graphs without electron loops were already processed

the 47 4-loop electron self-masses without electron loops



$$= \{23299 \text{ integrals}\} = \sum_{j=1}^{140} \frac{p_j(D)}{q_j(D)} M_j(D) \qquad 140 \text{ Master Integrals}$$

 $M_1$  M.I. with 11 denominators (still to be evaluated)  $p_1(D) =$ polynomial of degree 11  $q_1(D) = 5(D-1)(D-2)(D-3)(5D-16)(5D-18)(5D-22)$   $M_{140}$  M.I. with 4 denominators (factorizes into 4 1-loop tadpoles)  $p_{140}(D) =$ polynomial of degree 58

$$q_{140}(D) = 5184000D(D+4)(D+2)(D-1)^2(D-3)^5(D-4)^4$$
$$(D-5)^3(D-6)(D-8)(2D-5)^2(2D-7)^4(2D-9)^2(2D-11)$$
$$(2D-13)(3D-8)^2(3D-10)^3(3D-11)^2(3D-13)(4D-11)$$
$$(5D-12)(5D-13)(5D-14)(5D-16)(5D-17)(5D-18)$$
$$(5D-19)(5D-21)(5D-22)(7D-16)(10D^2-59D+86)$$

 $M_{140}$  is trivial!

$$M = M_{140} = \frac{1}{\epsilon^4 (1-\epsilon)^4}$$

with  $\epsilon = \frac{4-D}{2}$ .

The evaluation of the other MI's is not trivial.

A few analytic results are available, but the analytic evaluation of *all* the MI's does not seem to be at present a realistic possibility.

But a High precision numerical algorithm has been developped by S. Laporta a few yeas ago (2000).

Given a master integral M(D)

- choose a denominator in M(D)
- consider the set of the integrals M(n, D) where that denominator is raised to the power n
- construct a (system of) difference equation in n for the M(n, D)
- solve the system by expansion in factorial series or by Laplace transformation.

Advantages:

- No numerical integration
- The calculation of multi-loop integrals (in *any* number of loops) is reduced to sums of series in one variable.
- Arbitrary precision obtainable (such as 100-1000 digits).

as an example, consider  $M_{139}$ , which is *not* trivial; a semi-analytic approach gives

$$= M_{139} = -\frac{5}{2\epsilon^4} - \frac{45}{4\epsilon^3} - \frac{4255}{144\epsilon^2} - \frac{106147}{1728\epsilon}$$

$$+\frac{\pi\sqrt{3}}{240}\left(297C - 1477E\right) - \frac{2320981}{20736} + O(\epsilon)$$

$$C = \frac{4\pi^2}{27} \int_{0}^{1} \frac{dx}{\sqrt{1-x}} \, _2F_1^2\left(\begin{smallmatrix}\frac{1}{3}&\frac{2}{3}\\1\end{smallmatrix};x\right)$$

 $= 7.396\ 099\ 534\ 768\ 919\ 553\ 449\ 114\ 417\ 961\ 526\ 519\ 642\ldots$ 

$$E = \frac{4\pi^2}{27} \int_0^1 \frac{dx}{\sqrt{1-x}} {}_2F_1^2 \left( \frac{1}{3} - \frac{1}{3}; x \right)$$
  
= 2.376 887 326 184 666 003 152 855 958 761 330 926 023.

The Laporta High-precision numerical algorithm gives

$$= M_{139} = -\frac{5}{2\epsilon^4} - \frac{45}{4\epsilon^3} - \frac{4255}{144\epsilon^2} - \frac{106147}{1728\epsilon} - 141.72215618664768694996791$$

 $-521.14654568600250441775466\epsilon - 3347.9933650782886117865341\epsilon^2$ 

 $- 17951.3774774809944931097622\epsilon^3 - 101753.8165331173182139560386\epsilon^4 + \ldots$