

A Simple Subtraction Procedure for Calculation of the Anomalous Magnetic Moment of the Electron in QED

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▪ **Bogoliubov-Parasiuk theorem** [1956]

Removal of ultraviolet divergences by R-operation (defined by recurrence relations).

▪ **Forest formula:** V. Scherbina [1964], O. Zavyalov, B. Stepanov [1965], W. Zimmermann [1969]. “BPHZ renormalization”.

$$f^{\text{UV-free}} = (1 - K_1)(1 - K_2) \dots (1 - K_n) f$$

K_j transforms Feynman amplitude of j -th divergent subgraph (G_j) into its Taylor expansion up to $\omega(G_j)$ order at 0 (in momentum representation).

All terms with overlapping elements must be removed.

$$\omega(G) = \text{degree of UV divergence} = 4 - N_\mu - (3/2)N_e$$

UV-divergences are removed in Schwinger-parametric representation point-by-point, before integration, if $i\varepsilon$ in propagator denominators is fixed ($\varepsilon > 0$).

$$\text{Schwinger parameters: } 1/(x+i\varepsilon) = (1/i) \cdot \int_0^{+\infty} e^{i\alpha(x+i\varepsilon)} d\alpha$$

$\varepsilon \rightarrow 0 \Rightarrow$ IR-divergences.

AMM of the electron (theory and experiment)

The measured value [2011]:

$$a_e = 0.00115965218073(28)$$

The most accurate prediction (T. Kinoshita et al. [2015]):

$$a_e = a_e(QED) + a_e(hadronic) + a_e(electroweak),$$

$$a_e(QED) = \sum_{n \geq 1} \left(\frac{\alpha}{\pi} \right)^n a_e^{2n},$$

$$a_e^{2n} = A_1^{(2n)} + A_2^{(2n)}(m_e / m_\mu) + A_2^{(2n)}(m_e / m_\tau) + A_3^{(2n)}(m_e / m_\mu, m_e / m_\tau)$$

$$a_e = 0.001159652181643(25)(23)(16)(763)$$

($\alpha^{-1} = 137.035999049(90)$ – from experiments with rubidium atoms)

Uncertainties come from:

$$A_1^{(8)}, A_1^{(10)}, a_e(hadronic) + a_e(electroweak), \alpha$$

T. Aoyama, M. Hayakawa, T. Kinoshita, M. Nio, Tenth-Order Electron Anomalous Magnetic Moment – Contribution of Diagrams without Closed Lepton Loops, Physical Review D, 2015, V. 91, 033006.

My method was developed for computing $A_1^{(2n)}$

Universal QED contributions

$$a_e = a_e(QED) + a_e(hadroni) + a_e(electroweak),$$

$$a_e(QED) = \sum_{n \geq 1} \left(\frac{\alpha}{\pi} \right)^n a_e^{2n},$$

$$a_e^{2n} = A_1^{(2n)} + A_2^{(2n)}(m_e / m_\mu) + A_2^{(2n)}(m_e / m_\tau) + A_3^{(2n)}(m_e / m_\mu, m_e / m_\tau)$$

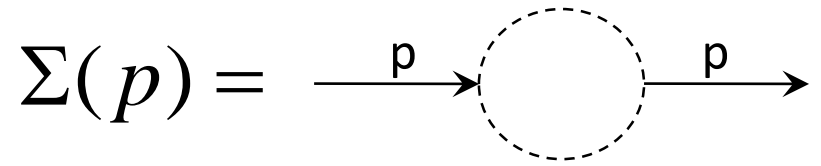
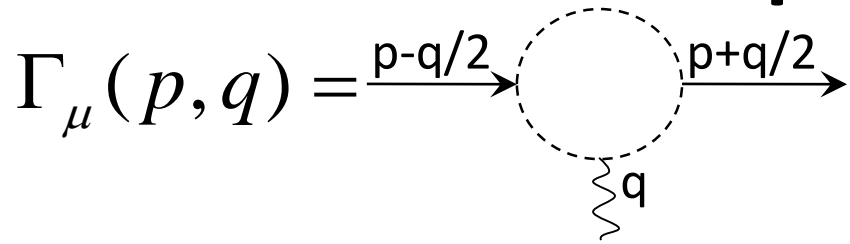
- J. Schwinger [1948], analytically: $A_1^{(2)} = 0.5$
- R. Karplus, N. Kroll [1949] – with a mistake
 - A. Petermann [1957], C. Sommerfield [1958], analytically:

$$A_1^{(4)} = -0.32847896579193..$$
- ~1965...~1975, 3 loops, numerically:
 1. M. Levine, J. Wright.
 2. R. Carroll, Y. Yao.
 3. T. Kinoshita, P. Cvitanović.
 - T. Kinoshita, P. Cvitanović [1974]: $A_1^{(6)} = 1.195 \pm 0.026$
- E. Remiddi, S. Laporta et al., ~1965...1996, analytically: $A_1^{(6)} = 1.181241456$
- T. Kinoshita et al., numerically, 1981...2015: $A_1^{(8)} = -1.9129884$
- T. Kinoshita et al., numerically, 2012...2015: $A_1^{(10)} = 7.795(336)$

The subtraction procedure

- FULLY AUTOMATED AT ANY ORDER OF PERTURBATION.
- UV and IR divergences are eliminated point-by-point in Feynman-parametric space for each individual Feynman diagram. No regularization is required.
- Subtraction by a forest formula with linear operators. Each operator transforms Feynman amplitude of some UV-divergent subdiagram G' (in momentum space) to the polynomial with the degree that is less or equal to $\omega(G')$.
- The subtraction is equivalent to the on-shell renormalization => no residual renormalizations, no calculations of renormalization constants, no other manipulations.

Operators



■ A – projector of AMM

$$\bar{u}_2 \Gamma_\mu(p, q) u_1 = \bar{u}_2 (f(q^2) \gamma_\mu - g(q^2) \sigma_{\mu\nu} q^\nu / (2m) + h(q^2) q_\mu) u_1$$

$$\sigma_{\mu\nu} = (\gamma_\mu \gamma_\nu - \gamma_\nu \gamma_\mu) / 2, \quad (p - q/2)^2 = (p + q/2)^2 = m^2$$

$$(\hat{p} - \hat{q}/2 - m) u_1 = (\hat{p} + \hat{q}/2 - m) u_2 = 0$$

$$A \Gamma_\mu = \gamma_\mu \lim_{q^2 \rightarrow 0} [g(q^2) + Ch(q^2)]$$

■ U – intermediate operator

$$\Gamma_\mu(p, 0) = a(p^2) \gamma_\mu + b(p^2) p_\mu + c(p^2) \hat{p} p_\mu + d(p^2) (\hat{p} \gamma_\mu - \gamma_\mu \hat{p}) \quad \Sigma(p) = r(p^2) + s(p^2) \hat{p}$$

$$U \Gamma_\mu = \gamma_\mu a(m^2)$$

$$U \Sigma = r(m^2) + s(m^2) \hat{p}$$

U preserves the Ward identity!

For other types of divergent subgraphs, U = Taylor expansion at 0 up to ω order.

■ L – on-shell renormalization for vertex-like subgraphs

$$L \Gamma_\mu = \gamma_\mu (a(m^2) + b(m^2) m + c(m^2) m^2)$$

Forest formula for AMM

A set of subgraphs of a diagram is called a **forest** if any two elements of this set don't overlap.

$F[G]$ – the set of all forests of UV-divergent subgraphs in G that contain G .

$\mathbb{I}[G]$ – the set of all vertex-like UV-divergent subgraphs in G that contains the vertex that is incident to the external photon line of G .

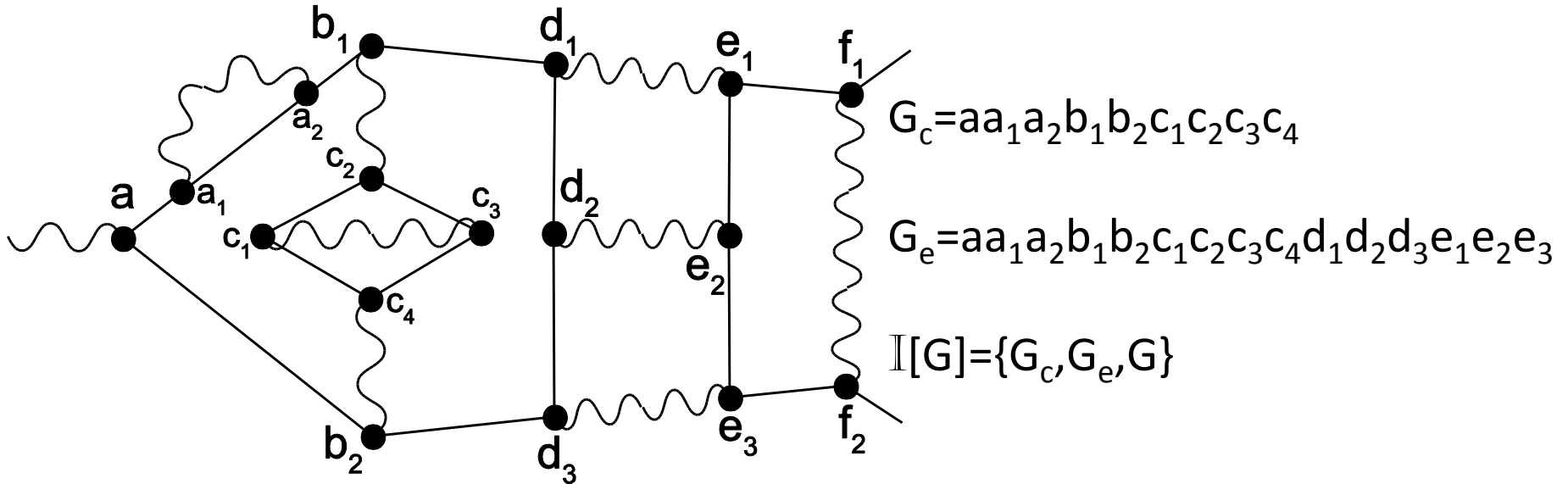
$$\tilde{f}_G = \sum_{\substack{F=\{G_1, \dots, G_n\} \in F[G] \\ G' \in \mathbb{I}[G] \cap F}} (-1)^{n-1} K_{G_1}^{F, G'} \dots K_{G_n}^{F, G'} f_G$$

$$K_{G''}^{F, G'} = \begin{cases} A_{G'} & \text{for } G' = G'' \\ U_{G''} & \text{for } G'' \notin \mathbb{I}[G] \text{ or } G'' \subseteq G', G'' \neq G' \\ L_{G''} & \text{for } G'' \in \mathbb{I}[G], G' \subseteq G'', G'' \neq G, G'' \neq G' \\ (L_{G''} - U_{G''}) & \text{for } G'' = G, G' \neq G \end{cases}$$

\bar{f}_G = coefficient before γ_μ in \tilde{f}_G

$$a_e = \sum_G \bar{f}_G$$

Example



Other UV-divergent subgraphs:

electron self-energy – a_1a_2 , vertex-like – $c_1c_2c_3, c_1c_3c_4$,

photon self-energy – $c_1c_2c_3c_4$,

photon-photon scattering – $G_d = aa_1a_2b_1b_2c_1c_2c_3c_4d_1d_2d_3$

$$\tilde{f}_G = \left[A_G (1 - U_{G_e}) (1 - U_{G_c}) - (L_G - U_G) A_{G_e} (1 - U_{G_c}) - (L_G - U_G) (1 - L_{G_e}) A_{G_c} \right] \cdot (1 - U_{G_d}) (1 - U_{c_1c_2c_3c_4}) (1 - U_{c_1c_2c_3} - U_{c_1c_3c_4}) (1 - U_{a_1a_2}) f_G$$

Feynman parameters

- Assign α_j to each line.
- New propagators

photon:
$$-\frac{g_{\mu\nu}}{p^2 + i\varepsilon} \Rightarrow i g_{\mu\nu} e^{i\alpha_j(p^2 + i\varepsilon)}$$

electron:
$$\frac{i(\hat{p} + m)}{p^2 - m^2 + i\varepsilon} \Rightarrow (\hat{p} + m) e^{i\alpha_j(p^2 - m^2 + i\varepsilon)}$$

- Apply the subtraction procedure with new propagators.
- Integrate (analytically) with respect to $\lambda = \alpha_1 + \dots + \alpha_n$.
- $\varepsilon \rightarrow 0$.
- Integrate (numerically) with respect to $\alpha_1, \dots, \alpha_n \geq 0$,
 $\alpha_1 + \dots + \alpha_n = 1$.

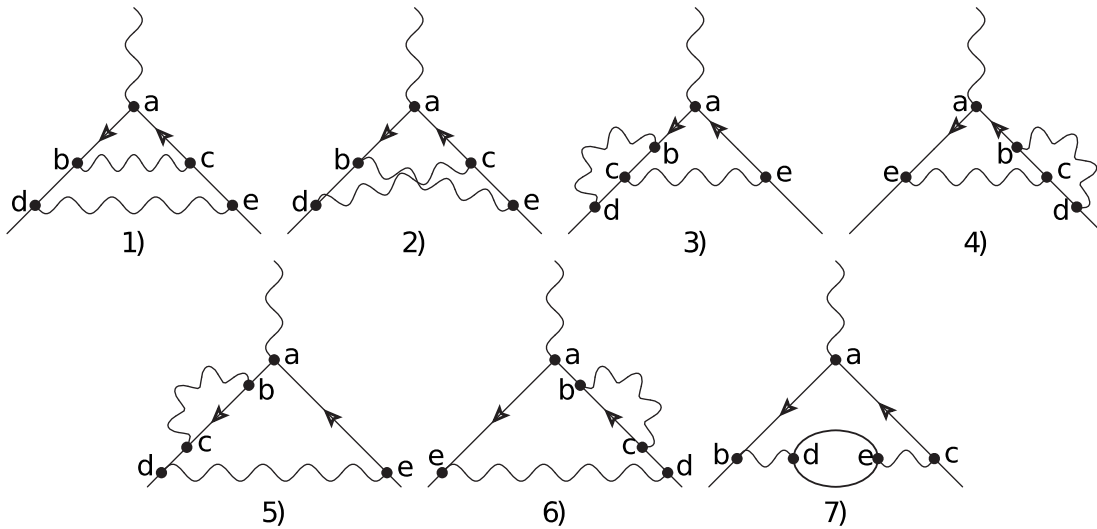
Realization

- 2 loops, 3 loops.
- Feynman gauge.
- D programming language – for the generator.
- C programming language – for the automatically generated code.
- Adaptive Monte-Carlo (homemade).
- 3 days of computation on a personal computer:

$$A_1^{(4)} = -0.32851(87)$$

$$A_1^{(6)} = 1.1802(85)$$

2-loop case

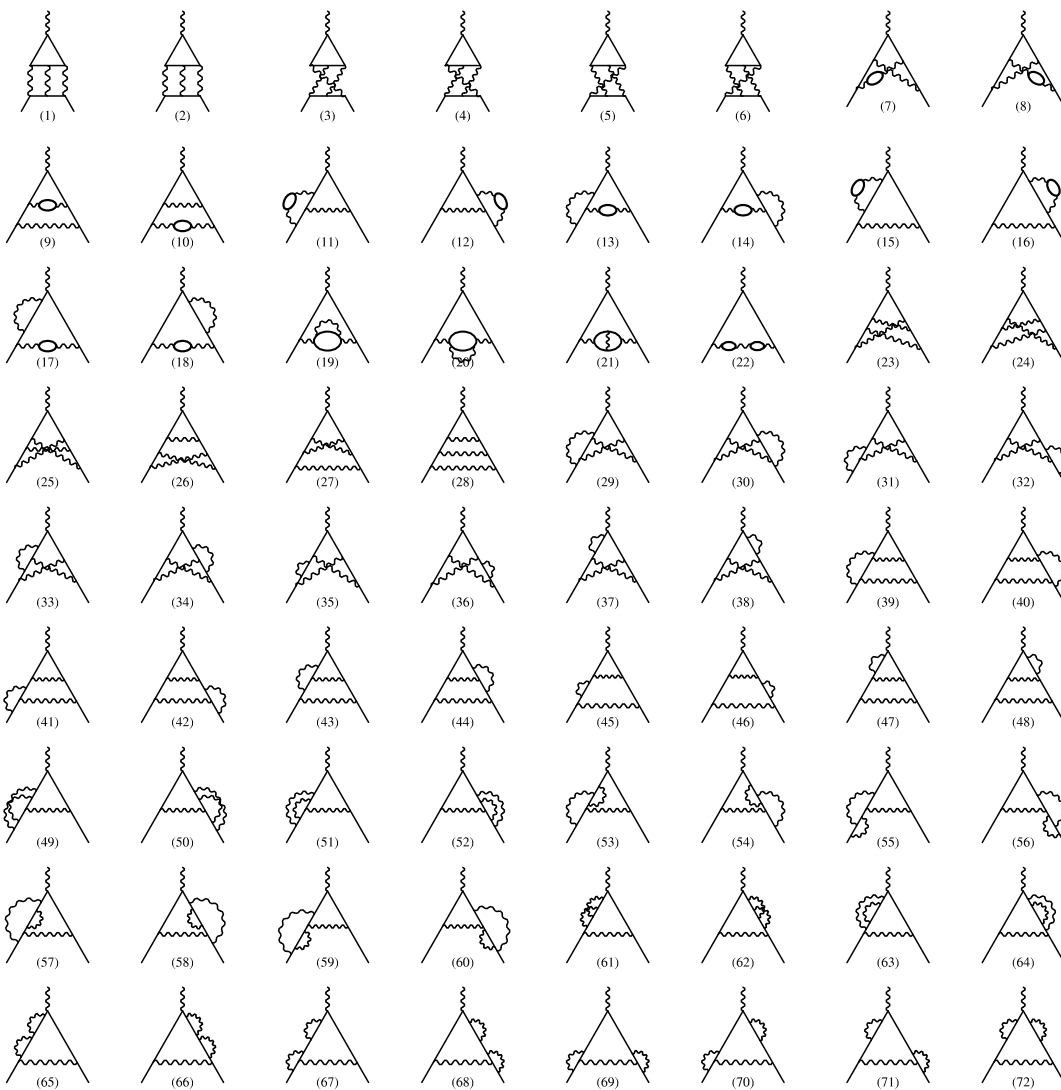


$$B\Sigma(p) = a(m^2) + mb(m^2) +$$

$$+ (\hat{p} - m)(b(m^2) + 2a'(m^2) + 2mb'(m^2)),$$

$$\Sigma(p) = a(p^2) + b(p^2)\hat{p}$$

#	My value	Analyt. value (Petermann, 1957)	Expression	On-shell renorm.	Difference
1	0.777455(52)	0.777478	$A_G - A_G U_{abc} - (L_G - U_G) A_{abc}$	$A_G - A_G L_{abc}$	$(L_G - U_G) A_{abc} - A_G (L_{abc} - U_{abc})$
2	-0.467626(44)	-0.467645	A_G	A_G	0
3	-0.032023(29)	-	$A_G - A_G U_{bcd}$	$A_G - A_G L_{bcd}$	$A_G (U_{abc} - L_{abc})$
4	-0.032033(29)	-	$A_G - A_G U_{bcd}$	$A_G - A_G L_{bcd}$	$A_G (U_{abc} - L_{abc})$
5	-0.294978(25)	-	$A_G - A_G U_{bc}$	$A_G - A_G B_{bc}$	$A_G (U_{bc} - B_{bc})$
6	-0.294998(24)	-	$A_G - A_G U_{bc}$	$A_G - A_G B_{bc}$	$A_G (U_{bc} - B_{bc})$
7	0.0156895(25)	0.0156874	$A_G - A_G U_{de}$	$A_G - A_G U_{de}$	0



3-loop Feynman diagrams for electron's AMM. Plot courtesy of F.Jegerlehner

Comparison with known analytical values

#	My value	Analyt. val.	Reference
1-6	0.3708(14)	0.3710	[10]
7-10	0.04989(20)	0.05015	[4,5]
11-12,15-16	-0.08782(15)	-0.08798	[2,4]
13-14,17-18	-0.11230(17)	-0.11234	[3,4]
19-21	0.05288(13)	0.05287	[1]
22	0.002548(20)	0.002559	[1]
23-24	1.8629(14)	1.8619	[11]
25	-0.02688(47)	-0.02680	[12]
26-27	-3.1764(22)	-3.1767	[8]
28	1.7888(19)	1.7903	[8]
29-30	-1.7579(10)	-1.7579	[12]
31-32,37-38	5.3559(27)	5.3576	[8,9]
33-34,37-38	0.4549(14)	0.4555	[8,11]
31-32,35-36	1.5436(34)	1.5416	[7,9]
33-36	-3.3573(24)	-3.3605	[7,11]
39-40	-0.33468(95)	-0.33470	[11]
41-48	-0.4030(41)	-0.4029	[6,7]
49-72	0.9529(53)	0.9541	[6-9,11,12]

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Thank you
for your attention!

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<http://arxiv.org/abs/1507.06435>