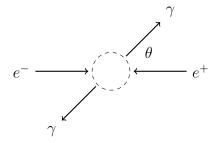
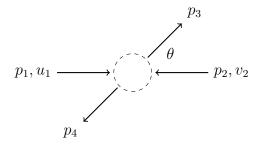
Electron positron annihilation creates two photons.



Here is the same diagram with momentum and spinor labels.



In a typical collider experiment the momentum vectors are

$$p_1 = \begin{pmatrix} E \\ 0 \\ 0 \\ p \end{pmatrix} \quad p_2 = \begin{pmatrix} E \\ 0 \\ 0 \\ -p \end{pmatrix} \quad p_3 = \begin{pmatrix} E \\ E \sin \theta \cos \phi \\ E \sin \theta \sin \phi \\ E \cos \theta \end{pmatrix} \quad p_4 = \begin{pmatrix} E \\ -E \sin \theta \cos \phi \\ -E \sin \theta \sin \phi \\ -E \cos \theta \end{pmatrix}$$
 inbound electron inbound positron outbound photon outbound photon

Symbol p is incident momentum, E is total energy $E = \sqrt{p^2 + m^2}$, and m is electron mass. Polar angle θ is the observed scattering angle. Azimuth angle ϕ cancels out in scattering calculations.

The spinors are

$$u_{11} = \begin{pmatrix} E + m \\ 0 \\ p \\ 0 \end{pmatrix} \qquad u_{12} = \begin{pmatrix} 0 \\ E + m \\ 0 \\ -p \end{pmatrix} \qquad v_{21} = \begin{pmatrix} -p \\ 0 \\ E + m \\ 0 \end{pmatrix} \qquad v_{22} = \begin{pmatrix} 0 \\ p \\ 0 \\ E + m \end{pmatrix}$$
inbound electron, spin up inbound electron, spin down inbound positron, spin up inbound positron, spin down inbound positron, spin up inbound positron, spin down inbound positron, spin down inbound positron, spin up inbound positron, spin down inbound positron, spin up inbound positron inbound

The spinors shown above are not individually normalized. Instead, a combined spinor normalization constant $N = (E + m)^2$ will be used.

The following formula computes a probability density $|\mathcal{M}_{ab}|^2$ for annihilation where a is the spin state of the inbound electron and b is the spin state of the inbound positron. The formula is from Feynman diagrams.

$$|\mathcal{M}_{ab}|^2 = \frac{e^4}{N} \left| -\frac{\bar{v}_{2b}\gamma^{\mu}(\not q_1 + m)\gamma^{\nu}u_{1a}}{t - m^2} - \frac{\bar{v}_{2b}\gamma^{\nu}(\not q_2 + m)\gamma^{\mu}u_{1a}}{u - m^2} \right|^2$$

Symbol e is electron charge and

$$q_1 = p_1 - p_3$$
$$q_2 = p_1 - p_4$$

Symbols t and u are Mandelstam variables

$$t = q_1^2 = (p_1 - p_3)^2$$

$$u = q_2^2 = (p_1 - p_4)^2$$

Let

$$a_1 = \bar{v}_{2b}\gamma^{\mu}(\not q_1 + m)\gamma^{\nu}u_{1a}$$
 $a_2 = \bar{v}_{2b}\gamma^{\nu}(\not q_2 + m)\gamma^{\mu}u_{1a}$

Then

$$|\mathcal{M}_{ab}|^2 = \frac{e^4}{N} \left| -\frac{a_1}{t - m^2} - \frac{a_2}{u - m^2} \right|^2$$

$$= \frac{e^4}{N} \left(-\frac{a_1}{t - m^2} - \frac{a_2}{u - m^2} \right) \left(-\frac{a_1}{t - m^2} - \frac{a_2}{u - m^2} \right)^*$$

$$= \frac{e^4}{N} \left(\frac{a_1 a_1^*}{(t - m^2)^2} + \frac{a_1 a_2^*}{(t - m^2)(u - m^2)} + \frac{a_1^* a_2}{(t - m^2)(u - m^2)} + \frac{a_2 a_2^*}{(u - m^2)^2} \right)$$

The expected probability density $\langle |\mathcal{M}|^2 \rangle$ is computed by summing $|\mathcal{M}_{ab}|^2$ over all spin and polarization states and then dividing by the number of inbound states. There are four inbound states. The sum over polarization states is already accomplished by contraction of aa^* over μ and ν .

$$\langle |\mathcal{M}|^2 \rangle = \frac{1}{4} \sum_{a=1}^2 \sum_{b=1}^2 |\mathcal{M}_{ab}|^2$$

$$= \frac{e^4}{4N} \sum_{a=1}^2 \sum_{b=1}^2 \left(\frac{a_1 a_1^*}{(t-m^2)^2} + \frac{a_1 a_2^*}{(t-m^2)(u-m^2)} + \frac{a_1^* a_2}{(t-m^2)(u-m^2)} + \frac{a_2 a_2^*}{(u-m^2)^2} \right)$$

Use the Casimir trick to replace sums over spins with matrix products.

$$f_{11} = \frac{1}{N} \sum_{a=1}^{2} \sum_{b=1}^{2} a_{1} a_{1}^{*} = \text{Tr}\left((\not p_{1} + m)\gamma^{\mu}(\not q_{1} + m)\gamma^{\nu}(\not p_{2} - m)\gamma_{\nu}(\not q_{1} + m)\gamma_{\mu}\right)$$

$$f_{12} = \frac{1}{N} \sum_{a=1}^{2} \sum_{b=1}^{2} a_{1} a_{2}^{*} = \text{Tr}\left((\not p_{1} + m)\gamma^{\mu}(\not q_{2} + m)\gamma^{\nu}(\not p_{2} - m)\gamma_{\mu}(\not q_{1} + m)\gamma_{\nu}\right)$$

$$f_{22} = \frac{1}{N} \sum_{a=1}^{2} \sum_{b=1}^{2} a_{2} a_{2}^{*} = \text{Tr}\left((\not p_{1} + m)\gamma^{\mu}(\not q_{2} + m)\gamma^{\nu}(\not p_{2} - m)\gamma_{\nu}(\not q_{2} + m)\gamma_{\mu}\right)$$

Hence

$$\langle |\mathcal{M}|^2 \rangle = \frac{e^4}{4} \left(\frac{f_{11}}{(t-m^2)^2} + \frac{f_{12}}{(t-m^2)(u-m^2)} + \frac{f_{12}^*}{(t-m^2)(u-m^2)} + \frac{f_{22}}{(u-m^2)^2} \right)$$

Run "annihilation-1.txt" to verify the Casimir trick for electron positron annihilation.

The following formulas are equivalent to the Casimir trick. (Recall that $a \cdot b = a^{\mu}g_{\mu\nu}b^{\nu}$)

$$f_{11} = 32(p_1 \cdot p_3)(p_1 \cdot p_4) - 32m^2(p_1 \cdot p_2) + 64m^2(p_1 \cdot p_3) + 32m^2(p_1 \cdot p_4) - 64m^4$$

$$f_{12} = 16m^2(p_1 \cdot p_3) + 16m^2(p_1 \cdot p_4) - 32m^4$$

$$f_{22} = 32(p_1 \cdot p_3)(p_1 \cdot p_4) - 32m^2(p_1 \cdot p_2) + 32m^2(p_1 \cdot p_3) + 64m^2(p_1 \cdot p_4) - 64m^4$$

In Mandelstam variables

$$s = (p_1 + p_2)^2$$

$$t = (p_1 - p_3)^2$$

$$u = (p_1 - p_4)^2$$

the formulas are

$$f_{11} = 8tu - 24tm^2 - 8um^2 - 8m^4$$

$$f_{12} = 8sm^2 - 32m^4$$

$$f_{22} = 8tu - 8tm^2 - 24um^2 - 8m^4$$

Run "annihilation-2.txt" to verify.

High energy approximation

When $E \gg m$ a useful approximation is to set m=0 and obtain

$$f_{11} = 8tu$$
$$f_{12} = 0$$
$$f_{22} = 8tu$$

Hence

$$\langle |\mathcal{M}|^2 \rangle = \frac{e^4}{4} \left(\frac{8tu}{t^2} + \frac{8tu}{u^2} \right)$$
$$= 2e^4 \left(\frac{u}{t} + \frac{t}{u} \right)$$

For m = 0 the Mandelstam variables are

$$s = 4E^{2}$$

$$t = -2E^{2}(1 - \cos \theta)$$

$$u = -2E^{2}(1 + \cos \theta)$$

Hence

$$\langle |\mathcal{M}|^2 \rangle = 2e^4 \left(\frac{1 + \cos \theta}{1 - \cos \theta} + \frac{1 - \cos \theta}{1 + \cos \theta} \right)$$

Cross section

The differential cross section is

$$\frac{d\sigma}{d\Omega} = \frac{\langle |\mathcal{M}|^2 \rangle}{64\pi^2 s} = \frac{e^4}{32\pi^2 s} \left(\frac{1 + \cos\theta}{1 - \cos\theta} + \frac{1 - \cos\theta}{1 + \cos\theta} \right), \quad s \gg m$$

Substituting $e^4 = 16\pi^2\alpha^2$ yields

$$\frac{d\sigma}{d\Omega} = \frac{\alpha^2}{2s} \left(\frac{1 + \cos\theta}{1 - \cos\theta} + \frac{1 - \cos\theta}{1 + \cos\theta} \right)$$

We can integrate $d\sigma$ to obtain a cumulative distribution function. Recall that

$$d\Omega = \sin\theta \, d\theta \, d\phi$$

Hence

$$d\sigma = \frac{\alpha^2}{2s} \left(\frac{1 + \cos \theta}{1 - \cos \theta} + \frac{1 - \cos \theta}{1 + \cos \theta} \right) \sin \theta \, d\theta \, d\phi$$

Let $I(\theta)$ be the following integral of $d\sigma$.

$$I(\theta) = \frac{2s}{2\pi\alpha^2} \int_0^{2\pi} \int d\sigma$$
$$= \int \left(\frac{1 + \cos\theta}{1 - \cos\theta} + \frac{1 - \cos\theta}{1 + \cos\theta}\right) \sin\theta \, d\theta, \quad a \le \theta \le \pi - a$$

Angular support is limited to an arbitrary a > 0 because I(0) and $I(\pi)$ are undefined. Assume that $I(\theta) - I(a)$ is computable given θ by either symbolic or numerical integration.

Let C be the normalization constant

$$C = I(\pi - a) - I(a)$$

Then the cumulative distribution function $F(\theta)$ is

$$F(\theta) = \frac{I(\theta) - I(a)}{C}, \quad a \le \theta \le \pi - a$$

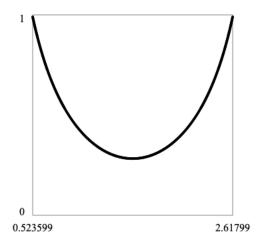
The probability of observing scattering events in the interval θ_1 to θ_2 can now be computed.

$$P(\theta_1 \le \theta \le \theta_2) = F(\theta_2) - F(\theta_1)$$

Probability density function $f(\theta)$ is the derivative of $F(\theta)$.

$$f(\theta) = \frac{dF(\theta)}{d\theta} = \frac{1}{C} \left(\frac{1 + \cos \theta}{1 - \cos \theta} + \frac{1 - \cos \theta}{1 + \cos \theta} \right) \sin \theta, \quad a \le \theta \le \pi - a$$

Run "annihilation-4.txt" to plot $f(\theta)$ for $a = \pi/6 = 30^{\circ}$.



Probability distribution for 30° bins $(a = 30^{\circ})$.

θ_1	θ_2	$P(\theta_1 \le \theta \le \theta_2)$
0°	30°	_
30°	60°	0.33
60°	90°	0.17
90°	120°	0.17
120°	150°	0.33
150°	180°	_

Data from DESY PETRA experiment

See www.hepdata.net/record/ins191231, Table 2, 14.0 GeV.

x	y
0.0502	0.09983
0.1505	0.10791
0.2509	0.12026
0.3512	0.13002
0.4516	0.17681
0.5521	0.1957
0.6526	0.279
0.7312	0.33204

Data x and y have the following relationship with the differential cross section formula.

$$x = \cos \theta \qquad y = \frac{d\sigma}{d\Omega}$$

To compute predicted values \hat{y} from the cross section formula, use $s = (14.0 \,\text{GeV})^2$. Multiply by $(\hbar c)^2$ to convert to SI and multiply by 10^{37} to convert square meters to nanobarns.

$$\hat{y} = \frac{\alpha^2}{2s} \left(\frac{1+x}{1-x} + \frac{1-x}{1+x} \right) \times (\hbar c)^2 \times 10^{37}$$

The following table shows predicted values \hat{y} .

x	y	\hat{y}
0.0502	0.09983	0.106325
0.1505	0.10791	0.110694
0.2509	0.12026	0.120005
0.3512	0.13002	0.135559
0.4516	0.17681	0.159996
0.5521	0.1957	0.198562
0.6526	0.279	0.262745
0.7312	0.33204	0.348884

The coefficient of determination R^2 measures how well predicted values fit the real data.

$$R^{2} = 1 - \frac{\sum (y - \hat{y})^{2}}{\sum (y - \bar{y})^{2}} = 0.98$$

The result indicates that the model $d\sigma$ explains 98% of the variance in the data.

Run "annihilation-3.txt" to verify.

Notes

Here are some notes on how the Eigenmath scripts work.

To convert a_1 and a_2 to Eigenmath code, it is instructive to write a_1 and a_2 in full component form.

$$a_1^{\mu\nu} = \bar{v}_{2\alpha}\gamma^{\mu\alpha}{}_\beta(\not\! q_1+m)^\beta{}_\rho\gamma^{\nu\rho}{}_\sigma u_1^\sigma \qquad a_2^{\nu\mu} = \bar{v}_{2\alpha}\gamma^{\nu\alpha}{}_\beta(\not\! q_2+m)^\beta{}_\rho\gamma^{\mu\rho}{}_\sigma u_1^\sigma$$

Transpose the γ tensors to form inner products over α and ρ .

$$a_1^{\mu\nu} = \bar{v}_{2\alpha}\gamma^{\alpha\mu}{}_\beta (\not q_1 + m)^\beta{}_\rho \gamma^{\rho\nu}{}_\sigma u_1^\sigma \qquad a_2^{\nu\mu} = \bar{v}_{2\alpha}\gamma^{\alpha\nu}{}_\beta (\not q_2 + m)^\beta{}_\rho \gamma^{\rho\mu}{}_\sigma u_1^\sigma$$

Convert transposed γ to Eigenmath code.

$$\gamma^{lpha\mu}{}_{eta} ~
ightarrow ~{
m gammaT}$$
 = transpose(gamma)

Then to compute a_1 we have

$$a_1 = \bar{v}_{2\alpha} \gamma^{\alpha\mu}{}_{\beta} (\not q_1 + m)^{\beta}{}_{\rho} \gamma^{\rho\nu}{}_{\sigma} u_1^{\sigma}$$

$$\rightarrow \quad \text{a1 = dot(v2bar[s2],gammaT,qslash1 + m I,gammaT,u1[s1])}$$

where s_1 and s_2 are spin indices. Similarly for a_2 we have

$$a_2 = \bar{v}_{2\alpha} \gamma^{\alpha\mu}{}_{\beta} (\rlap/q_2 + m)^{\beta}{}_{\rho} \gamma^{\rho\nu}{}_{\sigma} u_1^{\sigma}$$

$$\rightarrow \quad \text{a2 = dot(v2bar[s2],gammaT,qslash2 + m I,gammaT,u1[s1])}$$

In component notation the product $a_1a_1^*$ is

$$a_1 a_1^* = a_1^{\mu\nu} a_1^{*\mu\nu}$$

To sum over μ and ν it is necessary to lower indices with the metric tensor. Also, transpose a_1^* to form an inner product with ν .

$$a_1 a_1^* = a_1^{\mu\nu} a_{1\nu\mu}^*$$

Convert to Eigenmath code. The dot function sums over ν and the contract function sums over μ .

$$a_1 a_1^* \quad o \quad ext{all = contract(dot(al,gmunu,transpose(conj(al)),gmunu))}$$

Similarly for $a_2a_2^*$ we have

$$a_2 a_2^* \rightarrow \text{a22} = \text{contract(dot(a2,gmunu,transpose(conj(a2)),gmunu))}$$

The product $a_1 a_2^*$ does not require a transpose because $a_2 = a_2^{\nu\mu}$.

$$a_1^{\mu \nu} a_{2 \nu \mu}^* \quad o \quad {
m al2} \; = \; {
m contract(dot(a1,gmunu,conj(a2),gmunu))}$$

In component notation, a trace operator becomes a sum over an index, in this case α .

$$f_{11} = \operatorname{Tr}\left((\not p_1 + m) \gamma^{\mu} (\not q_1 + m) \gamma^{\nu} (\not p_2 - m) \gamma_{\nu} (\not q_1 + m) \gamma_{\mu} \right)$$
$$= (\not p_1 + m)^{\alpha}{}_{\beta} \gamma^{\mu\beta}{}_{\rho} (\not q_1 + m)^{\rho}{}_{\sigma} \gamma^{\nu\sigma}{}_{\tau} (\not p_2 - m)^{\tau}{}_{\delta} \gamma_{\nu}{}^{\delta}{}_{\eta} (\not q_1 + m)^{\eta}{}_{\xi} \gamma_{\mu}{}^{\xi}{}_{\alpha}$$

As before, transpose γ tensors to form inner products.

$$f_{11} = (\not\!p_1 + m)^\alpha{}_\beta \gamma^{\beta\mu}{}_\rho (\not\!q_1 + m)^\rho{}_\sigma \gamma^{\sigma\nu}{}_\tau (\not\!p_2 - m)^\tau{}_\delta \gamma^\delta{}_{\nu\eta} (\not\!q_1 + m)^\eta{}_\xi \gamma^\xi{}_{\mu\alpha}$$

This is the code for transposing γ .

$$\gamma^{\beta\mu}{}_{\beta} \rightarrow {
m gammaT = transpose(gamma)}$$
 $\gamma^{\delta}{}_{\nu\eta} \rightarrow {
m gammaL = transpose(dot(gmunu,gamma))}$

To convert f_{11} to Eigenmath code, use an intermediate variable T for the inner product.

$$T^{lpha\mu
u}{}_{
u\mulpha}$$
 $ightarrow$ T = dot(P1,gammaT,Q1,gammaT,P2,gammaL,Q1,gammaL)

Now sum over the indices of T. The innermost contract sums over ν then the next contract sums over μ . Finally the outermost contract sums over α .

$$f_{11}$$
 $ightarrow$ f11 = contract(contract(Contract(T,3,4),2,3))

Follow suit for f_{22} . For f_{12} the order of the rightmost μ and ν is reversed.

$$f_{12} = \operatorname{Tr}\left((\not p_1 + m)\gamma^{\mu}(\not q_2 + m)\gamma^{\nu}(\not p_2 - m)\gamma_{\mu}(\not q_1 + m)\gamma_{\nu}\right)$$

The resulting inner product is $T^{\alpha\mu\nu}_{\mu\nu\alpha}$ so the contraction is different.

$$f_{12} \rightarrow \text{f12} = \text{contract(contract(Contract(T,3,5),2,3))}$$

The innermost contract sums over ν followed by sum over μ then sum over α .