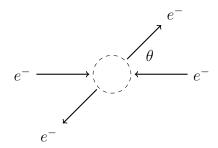
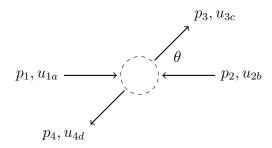
## Moller scattering

The following diagram shows the geometry of a Moller scattering experiment.



Here is the same diagram with momentum labels p and Dirac spinor labels u.



In center of mass coordinates the momentum vectors are

$$p_{1} = \begin{pmatrix} E \\ 0 \\ 0 \\ p \end{pmatrix} \qquad p_{2} = \begin{pmatrix} E \\ 0 \\ 0 \\ -p \end{pmatrix} \qquad p_{3} = \begin{pmatrix} E \\ p\sin\theta\cos\phi \\ p\sin\theta\sin\phi \\ p\cos\theta \end{pmatrix} \qquad p_{4} = \begin{pmatrix} E \\ -p\sin\theta\cos\phi \\ -p\sin\theta\sin\phi \\ -p\cos\theta \end{pmatrix}$$
electron 1 electron 2 electron 3

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

The spinors are

$$u_{11} = \begin{pmatrix} E+m \\ 0 \\ p \\ 0 \end{pmatrix} \qquad u_{21} = \begin{pmatrix} E+m \\ 0 \\ -p \\ 0 \end{pmatrix} \qquad u_{31} = \begin{pmatrix} E+m \\ 0 \\ p_{3z} \\ p_{3x}+ip_{3y} \end{pmatrix} \qquad u_{41} = \begin{pmatrix} E+m \\ 0 \\ p_{4z} \\ p_{4x}+ip_{4y} \end{pmatrix}$$

$$\stackrel{\text{electron 1}}{\underset{\text{spin up}}{\text{electron 3}}} \qquad u_{42} = \begin{pmatrix} 0 \\ E+m \\ 0 \\ p_{4z} \\ p_{4x}+ip_{4y} \end{pmatrix}$$

$$u_{12} = \begin{pmatrix} 0 \\ E+m \\ 0 \\ p \end{pmatrix} \qquad u_{22} = \begin{pmatrix} 0 \\ E+m \\ 0 \\ p \end{pmatrix} \qquad u_{32} = \begin{pmatrix} 0 \\ E+m \\ p_{3x}-ip_{3y} \\ -p_{3z} \end{pmatrix} \qquad u_{42} = \begin{pmatrix} 0 \\ E+m \\ p_{4x}-ip_{4y} \\ -p_{4z} \end{pmatrix}$$

$$\stackrel{\text{electron 1}}{\underset{\text{spin down}}{\text{electron 2}}} \qquad \stackrel{\text{electron 3}}{\underset{\text{spin down}}{\text{spin down}}} \qquad \stackrel{\text{electron 4}}{\underset{\text{spin down}}{\text{electron 4}}} \qquad \stackrel{\text{electron 4}}{\underset{\text{spin down}}{\text{spin down}}}$$

Spinor subscripts have 1 for spin up and 2 for spin down. The spinors are not individually normalized. Instead, a combined spinor normalization constant  $N = (E + m)^4$  will be used.

This is the probability density for spin state *abcd*. The formula is derived from Feynman diagrams for Moller scattering.

$$|\mathcal{M}_{abcd}|^2 = \frac{e^4}{N} \left| \frac{1}{t} (\bar{u}_{3c} \gamma^{\mu} u_{1a}) (\bar{u}_{4d} \gamma_{\mu} u_{2b}) - \frac{1}{u} (\bar{u}_{4d} \gamma^{\nu} u_{1a}) (\bar{u}_{3c} \gamma_{\nu} u_{2b}) \right|^2$$
no electron interchange electron interchange

Symbol e is electron charge. Symbols t and u are Mandelstam variables

$$t = (p_1 - p_3)^2$$
$$u = (p_1 - p_4)^2$$

Let

$$a_1 = (\bar{u}_{3c}\gamma^{\mu}u_{1a})(\bar{u}_{4d}\gamma_{\mu}u_{2b}), \quad a_2 = (\bar{u}_{4d}\gamma^{\nu}u_{1a})(\bar{u}_{3c}\gamma_{\nu}u_{2b})$$

Then

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

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

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

The expected probability density  $\langle |\mathcal{M}|^2 \rangle$  is computed by summing  $|\mathcal{M}_{abcd}|^2$  over all spin states and dividing by the number of inbound states. There are four inbound states.

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

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

The Casimir trick uses matrix arithmetic to compute sums.

$$f_{11} = \frac{1}{N} \sum_{abcd} a_1 a_1^* = \operatorname{Tr} \left( (\not p_3 + m) \gamma^{\mu} (\not p_1 + m) \gamma^{\nu} \right) \operatorname{Tr} \left( (\not p_4 + m) \gamma_{\mu} (\not p_2 + m) \gamma_{\nu} \right)$$

$$f_{12} = \frac{1}{N} \sum_{abcd} a_1 a_2^* = \operatorname{Tr} \left( (\not p_3 + m) \gamma^{\mu} (\not p_1 + m) \gamma^{\nu} (\not p_4 + m) \gamma_{\mu} (\not p_2 + m) \gamma_{\nu} \right)$$

$$f_{22} = \frac{1}{N} \sum_{abcd} a_2 a_2^* = \operatorname{Tr} \left( (\not p_4 + m) \gamma^{\mu} (\not p_1 + m) \gamma^{\nu} \right) \operatorname{Tr} \left( (\not p_3 + m) \gamma_{\mu} (\not p_2 + m) \gamma_{\nu} \right)$$

Hence

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

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

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

$$f_{12} = -32(p_1 \cdot p_2)^2 + 32m^2(p_1 \cdot p_2) + 32m^2(p_1 \cdot p_3) + 32m^2(p_1 \cdot p_4) - 32m^4$$

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

In Mandelstam variables

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

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

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

the formulas are

$$f_{11} = 8s^2 + 8u^2 - 64sm^2 - 64um^2 + 192m^4$$
  

$$f_{12} = -8s^2 + 64sm^2 - 96m^4$$
  

$$f_{22} = 8s^2 + 8t^2 - 64sm^2 - 64tm^2 + 192m^4$$

## High energy approximation

For high energy experiments  $E\gg m$  a useful approximation is to set m=0 and obtain

$$f_{11} = 8s^2 + 8u^2$$
$$f_{12} = -8s^2$$
$$f_{22} = 8s^2 + 8t^2$$

Hence

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

$$= \frac{e^4}{4} \left( \frac{8s^2 + 8u^2}{t^2} - \frac{-8s^2}{tu} - \frac{-8s^2}{tu} + \frac{8s^2 + 8t^2}{u^2} \right)$$

$$= 2e^4 \left( \frac{s^2 + u^2}{t^2} + \frac{2s^2}{tu} + \frac{s^2 + t^2}{u^2} \right)$$

Combine terms so  $\langle |\mathcal{M}|^2 \rangle$  has a common denominator.

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

For m = 0 the Mandelstam variables are

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

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

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

Hence

$$\langle |\mathcal{M}|^2 \rangle = 2e^4 \left( \frac{32E^8 \cos^4 \theta + 192E^8 \cos^2 \theta + 288E^8}{16E^8 (\cos \theta - 1)^2 (\cos \theta + 1)^2} \right)$$
$$= 4e^4 \frac{(\cos^2 \theta + 3)^2}{(\cos \theta - 1)^2 (\cos \theta + 1)^2}$$
$$= 4e^4 \frac{(\cos^2 \theta + 3)^2}{\sin^4 \theta}$$

The following equivalent formula can also be used.

$$\begin{split} \langle |\mathcal{M}|^2 \rangle &= 2e^4 \left( \frac{s^2 + u^2}{t^2} + \frac{2s^2}{tu} + \frac{s^2 + t^2}{u^2} \right) \\ &= 2e^4 \left( \begin{array}{c} \frac{1 + \cos^4(\theta/2)}{\sin^4(\theta/2)} &+ \frac{2}{\sin^2(\theta/2)\cos^2(\theta/2)} + \frac{1 + \sin^4(\theta/2)}{\cos^4(\theta/2)} \right) \\ &\text{no electron interchange} & \text{interaction term} & \text{electron interchange} \end{split}$$

## Cross section

The differential cross section is

$$\frac{d\sigma}{d\Omega} = \frac{\langle |\mathcal{M}|^2 \rangle}{4(4\pi\varepsilon_0)^2 s}$$

For the high energy approximation m = 0 we have

$$\langle |\mathcal{M}|^2 \rangle = 4e^4 \frac{(\cos^2 \theta + 3)^2}{\sin^4 \theta}$$
 and  $s = 4E^2$ 

Substitute for  $\langle |\mathcal{M}|^2 \rangle$ .

$$\frac{d\sigma}{d\Omega} = \frac{e^4}{(4\pi\varepsilon_0)^2 s} \frac{(\cos^2\theta + 3)^2}{\sin^4\theta}$$

Noting that

$$e^2 = 4\pi\varepsilon_0 \alpha \hbar c$$

we can also write

$$\frac{d\sigma}{d\Omega} = \frac{\alpha^2 (\hbar c)^2}{s} \frac{(\cos^2 \theta + 3)^2}{\sin^4 \theta}$$

We can integrate  $d\sigma$  to obtain a cumulative distribution function. Let  $I(\theta)$  be the following integral of  $d\sigma$ .

$$I(\theta) = \int \frac{(\cos^2 \theta + 3)^2}{\sin^4 \theta} \sin \theta \, d\theta$$

The result is

$$I(\theta) = -\frac{8\cos\theta}{\sin^2\theta} - \cos\theta$$

The cumulative distribution function is

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

Angular support is limited to an arbitrary a > 0 because I(0) and  $I(\pi)$  are undefined.

The probability of observing scattering events in the interval  $\theta_1$  to  $\theta_2$  is

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

## Notes

1. A. Zee page 134 has the cross section

$$\frac{d\sigma}{d\Omega} = \left(\frac{e^2}{4\pi}\right)^2 \frac{1}{8E^2} f(\theta)$$

where  $f(\theta)$  is the probability density function

$$f(\theta) = \frac{1 + \cos^4(\theta/2)}{\sin^4(\theta/2)} + \frac{2}{\sin^2(\theta/2)\cos^2(\theta/2)} + \frac{1 + \sin^4(\theta/2)}{\cos^4(\theta/2)}$$

The probability density function is equivalent to

$$f(\theta) = \frac{2(\cos^2 \theta + 3)^2}{\sin^4 \theta}$$

Hence for natural units  $\varepsilon_0 = \hbar = c = 1$  and  $e^2 = 4\pi\alpha$  the above cross section is equivalent to

$$\frac{d\sigma}{d\Omega} = \frac{\alpha^2 (\hbar c)^2}{4E^2} \frac{(\cos^2 \theta + 3)^2}{\sin^4 \theta}$$

2. In component notation, the trace operators of the Casimir trick become sums over a repeated index, in this case  $\alpha$ .

$$\begin{split} f_{11} &= \left( (\rlap{/}{p}_3 + m)^\alpha{}_\beta \gamma^{\mu\beta}{}_\rho (\rlap{/}{p}_1 + m)^\rho{}_\sigma \gamma^{\nu\sigma}{}_\alpha \right) \left( (\rlap{/}{p}_4 + m)^\alpha{}_\beta \gamma_\mu{}^\beta{}_\rho (\rlap{/}{p}_2 + m)^\rho{}_\sigma \gamma_\nu{}^\sigma{}_\alpha \right) \\ f_{12} &= (\rlap{/}{p}_3 + m)^\alpha{}_\beta \gamma^{\mu\beta}{}_\rho (\rlap{/}{p}_1 + m)^\rho{}_\sigma \gamma^{\nu\sigma}{}_\tau (\rlap{/}{p}_4 + m)^\tau{}_\delta \gamma_\mu{}^\delta{}_\eta (\rlap{/}{p}_2 + m)^\eta{}_\xi \gamma_\nu{}^\xi{}_\alpha \\ f_{22} &= \left( (\rlap{/}{p}_4 + m)^\alpha{}_\beta \gamma^{\mu\beta}{}_\rho (\rlap{/}{p}_1 + m)^\rho{}_\sigma \gamma^{\nu\sigma}{}_\alpha \right) \left( (\rlap{/}{p}_3 + m)^\alpha{}_\beta \gamma_\mu{}^\beta{}_\rho (\rlap{/}{p}_2 + m)^\rho{}_\sigma \gamma_\nu{}^\sigma{}_\alpha \right) \end{split}$$

To convert the above formulas to Eigenmath code, the  $\gamma$  tensors need to be transposed so that repeated indices are adjacent to each other. Also, multiply  $\gamma^{\mu}$  by the metric tensor to lower the index.

$$\gamma^{\beta\mu}_{\ \rho} \rightarrow {\rm gammaT = transpose(gamma)}$$
 $\gamma^{\beta}_{\ \mu\rho} \rightarrow {\rm gammaL = transpose(dot(gmunu,gamma))}$ 

Define the following  $4 \times 4$  matrices.

Then for  $f_{11}$  we have the following Eigenmath code. The contract function sums over  $\alpha$ .

$$(\not\!\!p_3 + m)^\alpha{}_\beta \gamma^{\mu\beta}{}_\rho (\not\!\!p_1 + m)^\rho{}_\sigma \gamma^{\nu\sigma}{}_\alpha \quad \rightarrow \quad \text{T1 = contract(dot(X3,gammaT,X1,gammaT),1,4)}$$
 
$$(\not\!\!p_4 + m)^\alpha{}_\beta \gamma_\mu{}^\beta{}_\rho (\not\!\!p_2 + m)^\rho{}_\sigma \gamma_\nu{}^\sigma{}_\alpha \quad \rightarrow \quad \text{T2 = contract(dot(X4,gammaL,X2,gammaL),1,4)}$$

Next, multiply then sum over repeated indices. The dot function sums over  $\nu$  then the contract function sums over  $\mu$ . The transpose makes the  $\nu$  indices adjacent as required by the dot function.

$$f_{11} = \operatorname{Tr}(\cdots \gamma^{\mu} \cdots \gamma^{\nu}) \operatorname{Tr}(\cdots \gamma_{\mu} \cdots \gamma_{\nu}) \quad o \quad \operatorname{contract(dot(T1, transpose(T2)))}$$

Follow suit for  $f_{22}$ .

$$(\not\!\!p_4 + m)^\alpha{}_\beta \gamma^{\mu\beta}{}_\rho (\not\!\!p_1 + m)^\rho{}_\sigma \gamma^{\nu\sigma}{}_\alpha \quad \rightarrow \quad \text{T1 = contract(dot(X4,gammaT,X1,gammaT),1,4)}$$
 
$$(\not\!\!p_3 + m)^\alpha{}_\beta \gamma_\mu{}^\beta{}_\rho (\not\!\!p_2 + m)^\rho{}_\sigma \gamma_\nu{}^\sigma{}_\alpha \quad \rightarrow \quad \text{T2 = contract(dot(X3,gammaL,X2,gammaL),1,4)}$$

Then

$$f_{22} = \operatorname{Tr}(\cdots \gamma^{\mu} \cdots \gamma^{\nu}) \operatorname{Tr}(\cdots \gamma_{\mu} \cdots \gamma_{\nu}) \quad \rightarrow \quad \operatorname{contract(dot(T1, transpose(T2)))}$$

The calculation of  $f_{12}$  begins with

$$\begin{split} (\not\!p_3 + m)^\alpha{}_\beta \gamma^{\mu\beta}{}_\rho (\not\!p_1 + m)^\rho{}_\sigma \gamma^{\nu\sigma}{}_\tau (\not\!p_4 + m)^\tau{}_\delta \gamma_\mu{}^\delta{}_\eta (\not\!p_2 + m)^\eta{}_\xi \gamma_\nu{}^\xi{}_\alpha \\ & \to \quad \text{T = contract(dot(X3,gammaT,X1,gammaT,X4,gammaL,X2,gammaL),1,6)} \end{split}$$

Then sum over repeated indices  $\mu$  and  $\nu$ .

$$f_{12} = \operatorname{Tr}(\cdots \gamma^{\mu} \cdots \gamma^{\nu} \cdots \gamma_{\mu} \cdots \gamma_{\nu}) \rightarrow \operatorname{contract(contract(T,1,3))}$$