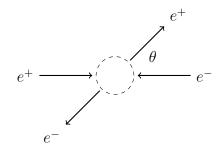
Bhabha scattering

Bhabha scattering is the interaction $e^- + e^+ \rightarrow e^- + e^+$.



In the center-of-mass frame we have the following momentum vectors where $E = \sqrt{p^2 + m^2}$.

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

Spinors for the inbound positron.

$$v_{11} = \frac{1}{\sqrt{E+m}} \begin{pmatrix} p \\ 0 \\ E+m \\ 0 \end{pmatrix} \qquad v_{12} = \frac{1}{\sqrt{E+m}} \begin{pmatrix} 0 \\ -p \\ 0 \\ E+m \end{pmatrix}$$
inbound e^+
spin up
inbound e^+
spin down

Spinors for the inbound electron.

$$u_{21} = \frac{1}{\sqrt{E+m}} \begin{pmatrix} E+m\\0\\-p\\0 \end{pmatrix} \qquad u_{22} = \frac{1}{\sqrt{E+m}} \begin{pmatrix} 0\\E+m\\0\\p \end{pmatrix}$$
inbound e^-
spin up
inbound e^-
spin down

Spinors for the outbound positron.

$$v_{31} = \frac{1}{\sqrt{E+m}} \begin{pmatrix} p_{3z} \\ p_{3x} + ip_{3y} \\ E+m \\ 0 \end{pmatrix} \qquad v_{32} = \frac{1}{\sqrt{E+m}} \begin{pmatrix} p_{3x} - ip_{3y} \\ -p_{3z} \\ 0 \\ E+m \end{pmatrix}$$
outbound e^+
spin up

Spinors for the outbound electron.

$$u_{41} = \frac{1}{\sqrt{E+m}} \begin{pmatrix} E+m\\0\\p_{4z}\\p_{4x}+ip_{4y} \end{pmatrix} \qquad u_{42} = \frac{1}{\sqrt{E+m}} \begin{pmatrix} 0\\E+m\\p_{4x}-ip_{4y}\\-p_{4z} \end{pmatrix}$$
outbound e^-
spin up
$$u_{41} = \frac{1}{\sqrt{E+m}} \begin{pmatrix} 0\\E+m\\p_{4x}-ip_{4y}\\-p_{4z} \end{pmatrix}$$
outbound e^-
spin down

The probability amplitude \mathcal{M}_{abcd} for spin state abcd is

$$\mathcal{M}_{abcd} = \mathcal{M}_{1abcd} + \mathcal{M}_{2abcd}$$

where

$$\mathcal{M}_{1abcd} = \frac{e^4}{s} (\bar{v}_{1a} \gamma^{\mu} u_{2b}) (\bar{u}_{4d} \gamma_{\mu} v_{3c}) \quad \mathcal{M}_{2abcd} = -\frac{e^4}{t} (\bar{v}_{1a} \gamma^{\nu} v_{3c}) (\bar{u}_{4d} \gamma_{\nu} u_{2b}),$$

Symbol e is elementary charge and

$$s = (p_1 + p_2)^2$$
$$t = (p_1 - p_3)^2$$

The expected probability density $\langle |\mathcal{M}|^2 \rangle$ is the average for all spin states.

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

Hence

$$\langle |\mathcal{M}|^2 \rangle = \frac{1}{4} \sum_{abcd} \left(\mathcal{M}_{1abcd} \mathcal{M}_{1abcd}^* + \mathcal{M}_{1abcd} \mathcal{M}_{2abcd}^* + \mathcal{M}_{2abcd} \mathcal{M}_{1abcd}^* + \mathcal{M}_{2abcd} \mathcal{M}_{2abcd}^* \right)$$

The Casimir trick uses matrix arithmetic to sum over spin states.

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

where

$$f_{11} = \operatorname{Tr}\left((\not p_1 - m)\gamma^{\mu}(\not p_2 + m)\gamma^{\nu}\right)\operatorname{Tr}\left((\not p_4 + m)\gamma_{\mu}(\not p_3 - m)\gamma_{\nu}\right)$$

$$f_{12} = \operatorname{Tr}\left((\not p_1 - m)\gamma^{\mu}(\not p_2 + m)\gamma^{\nu}(\not p_4 + m)\gamma_{\mu}(\not p_3 - m)\gamma_{\nu}\right)$$

$$f_{22} = \operatorname{Tr}\left((\not p_1 - m)\gamma^{\mu}(\not p_3 - m)\gamma^{\nu}\right)\operatorname{Tr}\left((\not p_4 + m)\gamma_{\mu}(\not p_2 + m)\gamma_{\nu}\right)$$

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

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

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

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

For high energy experiments such that $E \gg m$, let m = 0 and obtain

$$f_{11} = 32(p_1 \cdot p_3)^2 + 32(p_1 \cdot p_4)^2 = 64E^4(\cos^2\theta + 1)$$

$$f_{12} = -32(p_1 \cdot p_4)^2 = -32E^4(\cos\theta + 1)^2$$

$$f_{22} = 32(p_1 \cdot p_2)^2 + 32(p_1 \cdot p_4)^2 = 32E^4(\cos\theta + 1)^2 + 128E^4$$

For m = 0 the Mandelstam variables are

$$s = 4E^2$$
$$t = 2E^2(\cos\theta - 1)$$

Hence

$$\langle |\mathcal{M}|^2 \rangle = \frac{e^4}{4} \left(\frac{f_{11}}{s^2} - \frac{2f_{12}}{st} + \frac{f_{22}}{t^2} \right) = e^4 \left(\frac{\cos^2 \theta + 3}{\cos \theta - 1} \right)^2$$

Cross section

The differential cross section is

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

where

$$s = (p_1 + p_2)^2 = 4E^2$$

For high energy experiments we have

$$\langle |\mathcal{M}|^2 \rangle = e^4 \left(\frac{\cos^2 \theta + 3}{\cos \theta - 1} \right)^2$$

Hence for high energy experiments

$$\frac{d\sigma}{d\Omega} = \frac{e^4}{4(4\pi\varepsilon_0)^2 s} \left(\frac{\cos^2\theta + 3}{\cos\theta - 1}\right)^2$$

Noting that

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

we have

$$\frac{d\sigma}{d\Omega} = \frac{\alpha^2 (\hbar c)^2}{4s} \left(\frac{\cos^2 \theta + 3}{\cos \theta - 1} \right)^2$$

Noting that

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

we also have

$$d\sigma = \frac{\alpha^2 (\hbar c)^2}{4s} \left(\frac{\cos^2 \theta + 3}{\cos \theta - 1} \right)^2 \sin \theta \, d\theta \, d\phi$$

Let $S(\theta_1, \theta_2)$ be the following surface integral of $d\sigma$.

$$S(\theta_1, \theta_2) = \int_0^{2\pi} \int_{\theta_1}^{\theta_2} d\sigma$$

The solution is

$$S(\theta_1, \theta_2) = \frac{2\pi\alpha^2(\hbar c)^2}{4s} (I(\theta_2) - I(\theta_1))$$

where

$$I(\theta) = \frac{16}{\cos \theta - 1} - \frac{\cos^3 \theta}{3} - \cos^2 \theta - 9\cos \theta - 16\log(1 - \cos \theta)$$

The cumulative distribution function is

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

Angular support is reduced by an arbitrary angle a > 0 because I(0) is undefined.

The probability of observing scattering events in the interval θ_1 to θ_2 is

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

Let N be the total number of scattering events from an experiment. Then the number of scattering events in the interval θ_1 to θ_2 is predicted to be

$$NP(\theta_1 \le \theta \le \theta_2)$$

The probability density function is

$$f(\theta) = \frac{dF(\theta)}{d\theta} = \frac{1}{I(\pi) - I(a)} \left(\frac{\cos^2 \theta + 3}{\cos \theta - 1}\right)^2 \sin \theta$$

Data from SLAC SPEAR experiment

The following Bhabha scattering data is from SLAC-PUB-1501.

k	x_k	x_{k+1}	y
1	0.6	0.5	4432
2	0.5	0.4	2841
3	0.4	0.3	2045
4	0.3	0.2	1420
5	0.2	0.1	1136
6	0.1	0.0	852
7	0.0	-0.1	656
8	-0.1	-0.2	625
9	-0.2	-0.3	511
10	-0.3	-0.4	455
11	-0.4	-0.5	402
12	-0.5	-0.6	398

Column k is the bin number, column y is the number of scattering events, and

$$x_k = \cos \theta_k$$

The cumulative distribution function for this experiment is

$$F(\theta) = \frac{I(\theta) - I(\theta_1)}{I(\theta_{13}) - I(\theta_1)}$$

where

$$\theta_{13} = \arccos(-0.6), \quad \theta_1 = \arccos(0.6)$$

The scattering probability P_k is

$$P_k = F(\arccos(x_{k+1})) - F(\arccos(x_k))$$

Multiply P_k by total scattering events to obtain predicted number of events \hat{y}_k .

$$\sum y_k = 15773, \quad \hat{y}_k = 15773 \, P_k$$

The following table shows the predicted scattering events \hat{y} .

The coefficient of determination \mathbb{R}^2 measures how well predicted values fit the data.

$$R^{2} = 1 - \frac{\sum (y - \hat{y})^{2}}{\sum (y - \bar{y})^{2}} = 0.997$$

The result indicates that $F(\theta)$ explains 99.7% of the variance in the data.

Data from DESY PETRA experiment

See www.hepdata.net/record/ins191231, Table 3, 14.0 GeV.

x	y	
-0.7300	0.10115	
-0.6495	0.12235	
-0.5495	0.11258	
-0.4494	0.09968	
-0.3493	0.14749	
-0.2491	0.14017	
-0.1490	0.18190	
-0.0488	0.22964	
0.0514	0.25312	
0.1516	0.30998	
0.2520	0.40898	
0.3524	0.62695	
0.4529	0.91803	
0.5537	1.51743	
0.6548	2.56714	
0.7323	4.30279	

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

$$x = \cos \theta$$
, $y = \frac{d\sigma}{d\Omega}$ in units of nanobarns

The cross section formula is

$$\frac{d\sigma}{d\Omega} = \frac{\alpha^2}{4s} \left(\frac{\cos^2 \theta + 3}{\cos \theta - 1} \right)^2 \times (\hbar c)^2$$

To compute predicted values \hat{y} , multiply by 10^{37} to convert square meters to nanobarns.

$$\hat{y} = \frac{\alpha^2}{4s} \left(\frac{x^2 + 3}{x - 1} \right)^2 \times (\hbar c)^2 \times 10^{37}$$

The following table shows predicted values \hat{y} for $s = (14.0 \,\text{GeV})^2$.

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

$$R^{2} = 1 - \frac{\sum (y - \hat{y})^{2}}{\sum (y - \bar{y})^{2}} = 0.995$$

The result indicates that the model $d\sigma$ explains 99.5% of the variance in the data.

Notes

Here are a few notes about how the Eigenmath scripts work. In component notation the trace operators of the Casimir trick become sums over the repeated index α .

$$f_{11} = \left((\not p_1 - m)^{\alpha}{}_{\beta} \gamma^{\mu\beta}{}_{\rho} (\not p_3 - m)^{\rho}{}_{\sigma} \gamma^{\nu\sigma}{}_{\alpha} \right) \left((\not p_4 + m)^{\alpha}{}_{\beta} \gamma_{\mu}{}^{\beta}{}_{\rho} (\not p_2 + m)^{\rho}{}_{\sigma} \gamma_{\nu}{}^{\sigma}{}_{\alpha} \right)$$

$$f_{12} = (\not p_1 - m)^{\alpha}{}_{\beta} \gamma^{\mu\beta}{}_{\rho} (\not p_2 + m)^{\rho}{}_{\sigma} \gamma^{\nu\sigma}{}_{\tau} (\not p_4 + m)^{\tau}{}_{\delta} \gamma_{\mu}{}^{\delta}{}_{\eta} (\not p_3 - m)^{\eta}{}_{\xi} \gamma_{\nu}{}^{\xi}{}_{\alpha}$$

$$f_{22} = \left((\not p_1 - m)^{\alpha}{}_{\beta} \gamma^{\mu\beta}{}_{\rho} (\not p_2 + m)^{\rho}{}_{\sigma} \gamma^{\nu\sigma}{}_{\alpha} \right) \left((\not p_4 + m)^{\alpha}{}_{\beta} \gamma_{\mu}{}^{\beta}{}_{\rho} (\not p_3 - m)^{\rho}{}_{\sigma} \gamma_{\nu}{}^{\sigma}{}_{\alpha} \right)$$

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

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

Define the following 4×4 matrices.

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

$$(\not\!\!p_1 - m)^\alpha{}_\beta \gamma^{\mu\beta}{}_\rho (\not\!\!p_3 - m)^\rho{}_\sigma \gamma^{\nu\sigma}{}_\alpha \quad \rightarrow \quad \text{T1 = contract(dot(X1,gammaT,X3,gammaT),1,4)}$$

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

Next, multiply then sum over repeated indices. The dot function sums over ν then the contract function sums over μ . The transpose makes the ν indices adjacent as required by the dot function.

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

Follow suit for f_{22} .

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

Hence

$$f_{22} = \operatorname{Tr}(\cdots \gamma^{\mu} \cdots \gamma^{\nu}) \operatorname{Tr}(\cdots \gamma_{\mu} \cdots \gamma_{\nu}) \quad \rightarrow \quad \texttt{f22 = contract(dot(T1,transpose(T2)))}$$

The calculation of f_{12} begins with

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

Then sum over repeated indices μ and ν .

$$f_{12} = \operatorname{Tr}(\cdots \gamma^{\mu} \cdots \gamma^{\nu} \cdots \gamma_{\mu} \cdots \gamma_{\nu}) \quad o \quad ext{f12} = ext{contract(Contract(T,1,3))}$$