



Weak Decays of Doubly Heavy Baryons to Singly Heavy Baryons with FCNC

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In this work, we are examining the following decays:

- $\Xi_{ccq} \rightarrow \Sigma_{cuq}^*$
- $\Xi_{bbq} \rightarrow \Sigma_{bdq}^*$
- $\Xi_{bbq} \rightarrow \Xi_{bsq}^*$
- $\Omega_{bbs} \rightarrow \Omega_{bss}^*$
- $\Omega_{bbs} \rightarrow \Xi_{bds}^*$
- $\Omega_{ccs} \rightarrow \Xi_{cus}^*$

- Decays induced by flavor-changing neutral current take place at the loop level.
- All the QCD dynamics are encoded in the hadronic matrix elements induced by the weak current.
- These matrix elements are parameterized in terms of the form factors.
- Determining these form factors constitutes a central problem in studying such a class of decays.

- Matrix elements are defined as

$$\langle B_Q l^+ l^- | \mathcal{H}_{eff} | B_{QQ} \rangle$$

- where \mathcal{H}_{eff} for $b \rightarrow s, d$ is

$$\mathcal{H}_{eff} = \frac{G_F}{\sqrt{2}} V_{tb} V_{tq}^* \sum_i C_i^b(\mu) \mathcal{O}_i^b(\mu)$$

and for $c \rightarrow u$ is

$$\mathcal{H}_{eff} = \frac{G_F}{\sqrt{2}} \sum_{j=b,s,d} V_{cj} V_{uj}^* \sum_i C_i^c(\mu) \mathcal{O}_i^c(\mu)$$

- To calculate the matrix element, we need to calculate

$$\langle B_Q | \bar{q} \Gamma_\mu Q | B_{QQ} \rangle$$

where $\Gamma_\mu = \sigma_{\mu\nu}(1 - \gamma_5)q^\nu$ or $\Gamma_\mu = \gamma_\mu(1 - \gamma_5)$

$$\begin{aligned} \langle B_Q(p) | \bar{q} \gamma_\mu (\gamma_5) | B_{QQ}(p+q) \rangle = \\ \bar{u}^\alpha(p) \left[(q_\alpha p_\mu - p q g_{\alpha\mu} F_1^{V(A)} + (q_\alpha \gamma_\mu - \not{q} g_{\alpha\mu}) F_2^{V(A)} \right. \\ \left. + (q_\alpha q_\mu - q^2 g_{\alpha\mu} F_3^{V(A)} + q_\alpha q_\mu F_0^{V(A)}) \right] (\gamma_5) u(p+q) \end{aligned}$$

$$\begin{aligned} \langle B_Q(p) | i\bar{q}\sigma^{\mu\nu} Q | B_{QQ}(p+q) \rangle = \\ i\bar{u}_\alpha \left((p+q)^\alpha (X_1(-i\sigma^{\mu\nu}) + X_2\gamma^{[\mu}(p+q)^{\nu]}) \right. \\ \left. + X_3\gamma^{[\mu}p^{\nu]} + X_4(p+q)^{[\mu}p^{\nu]} \right) \\ \left. + X_5g^{\alpha[\mu}\gamma^{\nu]} + X_6g^{\alpha[\mu}(p+q)^{\nu]} + X_7g^{\alpha[\mu}p^{\nu]} \right) \end{aligned}$$

(Hiller and Zwicky, 2021)

$$T_1^V = 2X_1 + \Delta_-(X_2 + X_3) + q^2 X_4 + X_5 + X_6$$

$$T_2^V = -\Delta_+ X_1 + \frac{1}{2}(m_2^2 - q^2 - 1)X_2 + \frac{1}{2}(m_2^2 + q^2 - 1)X_3 + X_5$$

$$T_3^V = -X_1 + \Delta_- X_2 - \frac{1}{2}(m_2^2 + q^2 - 1)X_4 - X_6$$

$$T_1^A = 2X_A + \Delta_-(X_2 + X_3) + X_6 + X_7$$

$$T_2^A = -\Delta_- X_1 + \frac{1}{2}(q^2 + 2m_2^2 - m^2 - 1)(X_2 + X_3) + X_5 + m_2(X_6 + X_7)$$

$$T_3^A = -X_1 + m_2(X_2 + X_3)$$

where $\Delta_{\pm} = m_1 \pm m_2$ and m_1 is the mass the mass of the B_{QQ} and m_2 is the mass of the B_Q

- Collecting all, the transition amplitude can be written as

$$\mathcal{M} = \bar{u}_\alpha(p) \left\{ (\gamma_5)(\mathcal{F}_i^{A(V)} \mathcal{P}_i^{\alpha\mu} + C_9 \mathcal{P}_0^{\alpha\mu} F_0^{A(V)}) \bar{\ell} \gamma_\mu \ell \right. \\ \left. + (\gamma_5)(\mathcal{F}_4^{A(V)} \mathcal{P}_1^{\alpha\mu} + \mathcal{F}_5^{A(V)} \mathcal{P}_2^{\alpha\mu} \right. \\ \left. + \mathcal{F}_6^{A(V)} \mathcal{P}_3^{\alpha\mu} + C_{10} \mathcal{P}_0^{\alpha\mu} F_0^{A(V)}) \bar{\ell} \gamma_\mu \gamma_5 \ell \right\}$$

where $\mathcal{F}_i^{V(A)} = (-)C_9 F_i^{V(A)} - \frac{2m_Q}{q^2} C_7 T_i^{V(A)}$ with $i = 1, 2, 3$ and $\mathcal{F}_{4,5,6}^{V(A)} = (-)C_{10} F_{1,2,3}^{V(A)}$

- Using the sum rules method to calculate the form factors, we need the correlation function

$$\Pi_\mu = i \int d^4x e^{iqx} \langle B_Q(p) | \mathcal{T} \left\{ \bar{q}(x) \Gamma_\mu Q(x) j_{QQ}^\dagger(0) \right\} | 0 \rangle$$

where $j_{QQ} = \varepsilon^{abc} (Q^{aT} C \gamma^\alpha Q^b) \gamma_\alpha \gamma_5$ (Aliev et al., 2022)

- Using Wick theorem to calculate the correlation function with the QCD parameters, we contract the $Q(x)$ and the $\bar{Q}(0)$ from the $j_{QQ}^\dagger(0)$ we are left with matrix elements

$$\langle B_Q | \bar{q}_\alpha(x) \bar{q}'_\beta(0) \bar{Q}_\gamma(0) | 0 \rangle$$

where α, β and γ are spinor indices. These matrix elements are calculated in (Ali et al., 2013)

- We can also calculate the correlation function in terms of hadronic transition matrix elements. To calculate, we need to correlate the expression of the correlation function in the region $(p + q)^2 > 0$ and $(p + q)^2 \ll 0$ using the spectral representation.

$$\Pi((p + q)^2) = \int_0^\infty ds \frac{\rho(s, q^2, p^2)}{s - (p + q)^2} + \text{polynomial in } (p + q)^2$$

where $\rho(s, q^2, p^2)$ is the spectral density.

$$\rho^{phen}(s, q^2, p^2) = \langle B_Q | \bar{q} \Gamma_\mu Q | B_{QQ} \rangle \langle B_Q Q | j_{QQ}^\dagger | 0 \rangle \delta(s - m_{QQ}^2) + \rho^{higher}(s)$$

where $\rho^{higher}(s)$ represents the higher state and continuum contributions.

- To eliminate the polynomials and higher state and continuum contributions, Borel transformation and continuum subtraction can be applied. Finally, we get

$$\int_0^{s_0} ds e^{-s/M^2} \rho^{QCD}(s, q^2, p^2) = \int_0^{s_0} ds e^{-s/M^2} \rho^{phen}(s, q^2, p^2)$$

- To calculate the ρ^{phen} we need $\langle B_{QQ} | j_{QQ}^\dagger | 0 \rangle$ and it can be calculated with the following correlation function

$$\Pi = i \int d^4x e^{ipx} \langle 0 | \mathcal{T} \{ j_{QQ}(x) j_{QQ}^\dagger(0) \} | 0 \rangle$$

- After calculating both sides, we get different structures consisting of p and q and gamma matrices (i.e., $\gamma_\mu \not{q} q_\kappa$).
- Equating the coefficients of the structures on both sides, we can finally find the numerical values of the form factors.

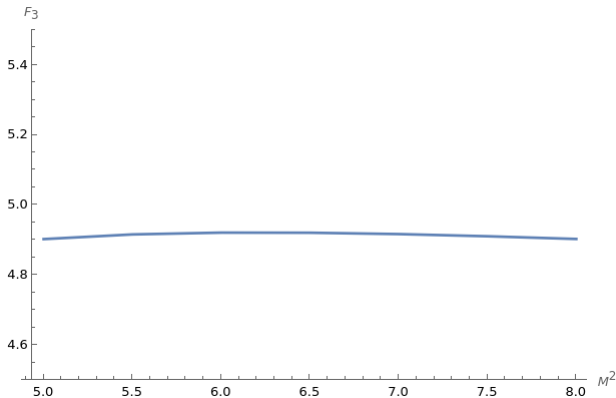


Figure: M^2 dependence

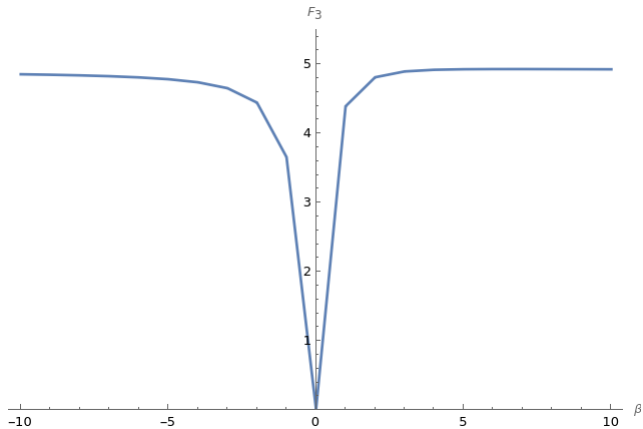


Figure: β dependency

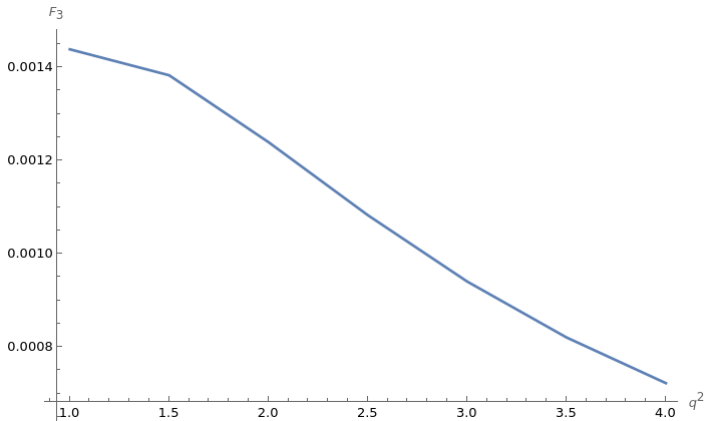


Figure: q^2 dependency

- We need the q^2 dependency of the form factors up to $(m_{QQ} - m_Q)^2$.
- Using the numerical values of the form factors for different q^2 values, we will find a fit

$$F(q^2) = \frac{F(0)}{1 - \frac{q^2}{m_{fit}^2} - \delta \left(\frac{q^2}{m_{fit}^2} \right)^2}$$

where m_{fit} and δ are fit parameters.

- Using these results, we will calculate the matrix elements.

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- Hiller, G., & Zwicky, R. (2021). Endpoint relations for baryons. *JHEP*, *11*, 073. [https://doi.org/10.1007/JHEP11\(2021\)073](https://doi.org/10.1007/JHEP11(2021)073)



Thank You
for your attention.

Do you have any question?