A New Approach for Delta Form Factors

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Calculating Form Factors

Difficulties in lattice calculations: Finite Volume, disconnected diagrams... Extracting ground states

Generalized Pencil-of-Function Method



 $\langle \Delta(p') | V^{\mu} | \Delta(p) \rangle = \bar{u}_{\alpha}(p') \Gamma^{\alpha \mu \beta} u_{\beta}(p)$



$$\Gamma^{\alpha\beta\mu}_{\gamma\Delta\Delta}(p',p) = -e \left\{ e_{\Delta} F_{1}^{*}(Q^{2}) g^{\alpha\beta} \gamma^{\mu} \right.$$

$$+ \frac{i}{2M_{\Delta}} \left[F_{2}^{*}(Q^{2}) g^{\alpha\beta} + F_{4}^{*}(Q^{2}) \frac{q^{\alpha}q^{\beta}}{(2M_{\Delta})^{2}} \right] \sigma^{\mu\nu} q_{\nu}$$

$$+ \frac{F_{3}^{*}(Q^{2})}{(2M_{\Delta})^{2}} \left[q^{\alpha}q^{\beta}\gamma^{\mu} - \frac{1}{2}q \cdot \gamma \left(g^{\alpha\mu}q^{\beta} + g^{\beta\mu}q^{\alpha} \right) \right] \right\}$$

At $Q^2 = 0$.



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Magnetic Dipole



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Charge Electric Quadrupole Magnetic Dipole



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Charge Electric Quadrupole Magnetic Dipole

Magnetic Octupole

From the PDG

$\mu_{\Delta^{++}} = (5.6 \pm 1.9)\mu_N$

 $\mu_{\Delta^+} = (2.7 \pm 3.5)\mu_N$

Even getting the magnetic moment experimentally is difficult for the Delta

The lattice is essential in determining the form factors in the absence of experimental data

On the lattice, we calculate n-point correlators:

 $C^{3pt}(t_i, t, t_f, \mathbf{p}_i, \mathbf{p}_f) = FT[\langle 0 | \chi(t_f) J_{\mu}(t) \chi(t_i) | 0 \rangle]$ For $t_f \gg t \gg t_i$: $C^{3pt}(t_i, t, t_f, \mathbf{p}_i, \mathbf{p}_f) \rightarrow Z(\mathbf{p}_i, \mathbf{p}_f) e^{-E_f(t_f - t)} e^{-E_i(t - t_i)}$ $\times \langle \Delta(p_f) | J_{\mu}(0) | \Delta(p_i) \rangle$ On the lattice, we calculate n-point correlators:

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And Z includes overlap factors that also arise in the two-point functions

 $C^{2pt}(t_i, t_f, \mathbf{p}) \to A'e^{-(t_f - t)E}$

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proton

<u>2</u>m $G_E(p^2) = F_1(p^2)$ F_{2}

 $G_E(p^2=0)$



Excited state contamination



Higher momenta





 $\frac{1}{Z_V} = 1.05(1)$

Not unity because we're using the local current

Errors are highly underestimated



The problem: excited state contamination

Are we at large enough time separation to only see ground state?

Yes for zero momentum but questionable for higher momenta...

Obligatory slide of lattice numbers given in unphysical units

Anisotropic Clover Lattices (via JLab)

$$\frac{a_s}{a_t} \approx 3.5 \qquad a_t^{-1} \approx 5.5 \text{ GeV}$$

Volumes: $(16^3, 20^3, 24^3) \times 128$ $390 \text{ (MeV)} \le m_{\pi}$ $m_{\Lambda} \approx 1.4 - 1.5 \text{ GeV}$

Anisotropy allows for better resolution in time (Good for excited states and baryons)

The problem also shows itself in "effective mass" plots



Excited state contamination

There are many operators that can be used

The only requirement is that they have a good overlap with the state of interest There are many operators that can be used

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One approach is to enumerate a large number of operators, and calculate a matrix of correlators

 $C_{ij}(t) = \langle 0|O_i(t)O_j(0)|0\rangle$

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Solve the Generalized Eigenvalue Problem

 $C(t)x = \lambda(t)C(t_0)x$

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One can show that the eigenvalues behave like

$$\lambda_i(t,t_0) \sim e^{-m_i(t-t_0)} + \cdots$$

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$$\lambda_i(t,t_0) \sim e^{-m_i(t-t_0)} + \cdots$$

 t_0 should be chosen such that $C(t_0)$ contains (ideally) all of the states in the correlator, no more, no less



Ground state has same behavior

Generalized PoF Method

Hua, Sakar (1989) Sarkar, Pereira (1995) **Generalized PoF Method**

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then so is

 $O_{\Delta}^{\tau}(t) \equiv e^{H\tau} O_{\Delta}(t) e^{-H\tau}$

Using this, we can consider

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Using this, we can consider $O^{\tau}_{\Delta}(t), O_{\Delta}(t)$ distinct operators and make a matrix of correlators from a single two-point function: $C(t) = \langle 0 | O_{\Delta}(t) O_{\Delta}^{\dagger}(0) | 0 \rangle$ $C(t+\tau) = \langle 0 | O^{\tau}_{\Lambda}(t) O^{\dagger}_{\Lambda}(0) | 0 \rangle$ $= \langle 0 | O_{\Delta}(t+\tau) O_{\Delta}^{\dagger}(0) | 0 \rangle$ $C(t+2\tau) = \langle 0|O^{\tau}_{\Lambda}(t)(O^{\tau}_{\Lambda})^{\dagger}(0)|0\rangle$ $= \langle 0 | O_{\Delta}(t+2\tau) O_{\Delta}^{\dagger}(0) | 0 \rangle$

In fact, we could do this as much as we want

 $\begin{pmatrix} C(t) & C(t+\tau) & C(t+2\tau) & \cdots & C(t+n\tau) \\ C(t+\tau) & C(t+2\tau) & C(t+3\tau) & \cdots & C(t+(n+1)\tau) \\ \end{pmatrix} \\ C(t+2\tau) & C(t+3\tau) & C(t+4\tau) & \cdots & \vdots \\ \vdots & \vdots & \ddots & \vdots \\ C(t+n\tau) & C(t+(n+1)\tau) & C(t+(n+2)\tau) & \cdots & C(t+2n\tau) \end{pmatrix}$

Where *n* is the number of shifts we perform and τ is the amount by which we shift

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Where *n* is the number of shifts we perform and τ is the amount by which we shift

In principle, we could choose $n=T/2, \ au=1$ with T the lattice time length Large correlations and noise and linear

dependence make this a bad idea



n, au m |Q|fit range 0.2770(50) [0.78] (29,40) 0, 00.2825(17) [0.84] (5,28) 1, 4 $2,4 \quad 0.2823(17) \quad [0.87] \quad (5,28)$ 2,2 0.2838(15) [0.74] (5,30) 0.2799(29)[0.90](5,22)2, 8Can we do better with more operators?

Not with only local operators...

As for the three-point correlators

$$\mathbf{C}^{3pt}(t_i, t, t_f) = \begin{pmatrix} C^{3pt}(t_i, t, t_f) & C^{3pt}(t_i, t, t_f + \tau) \\ C^{3pt}(t_i, t + \tau, t_f + \tau) & C^{3pt}(t_i, t + \tau, t_f + 2\tau) \end{pmatrix}$$

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Once the two-point correlator matrix is diagonalized with vectors V:

 $\mathbf{C}_{\text{diag}}^{3pt}(t_i, t, t_f) = V^{-1} \mathbf{C}^{3pt}(t_i, t, t_f) V$

Conclusions

Using Pencil-of-Function techniques is much better for getting the ground state than using multiple operators

We have a good determination of $G_E(q^2)$, and need to extract other form factors (and do so on larger volumes)

Hopefully applying GPoF techniques to the threepoint correlators will improve the signal even more (and systematics)