

Resonances and Lattice QCD

David Richards (Jefferson Laboratory/LHPC)

- Motivation
- Review
- Recipe
 - Variational Method
 - Group-theoretical methods
 - Numerical implementation
- Exploratory Tests
- Conclusions and Future Plans



LHP Collaboration

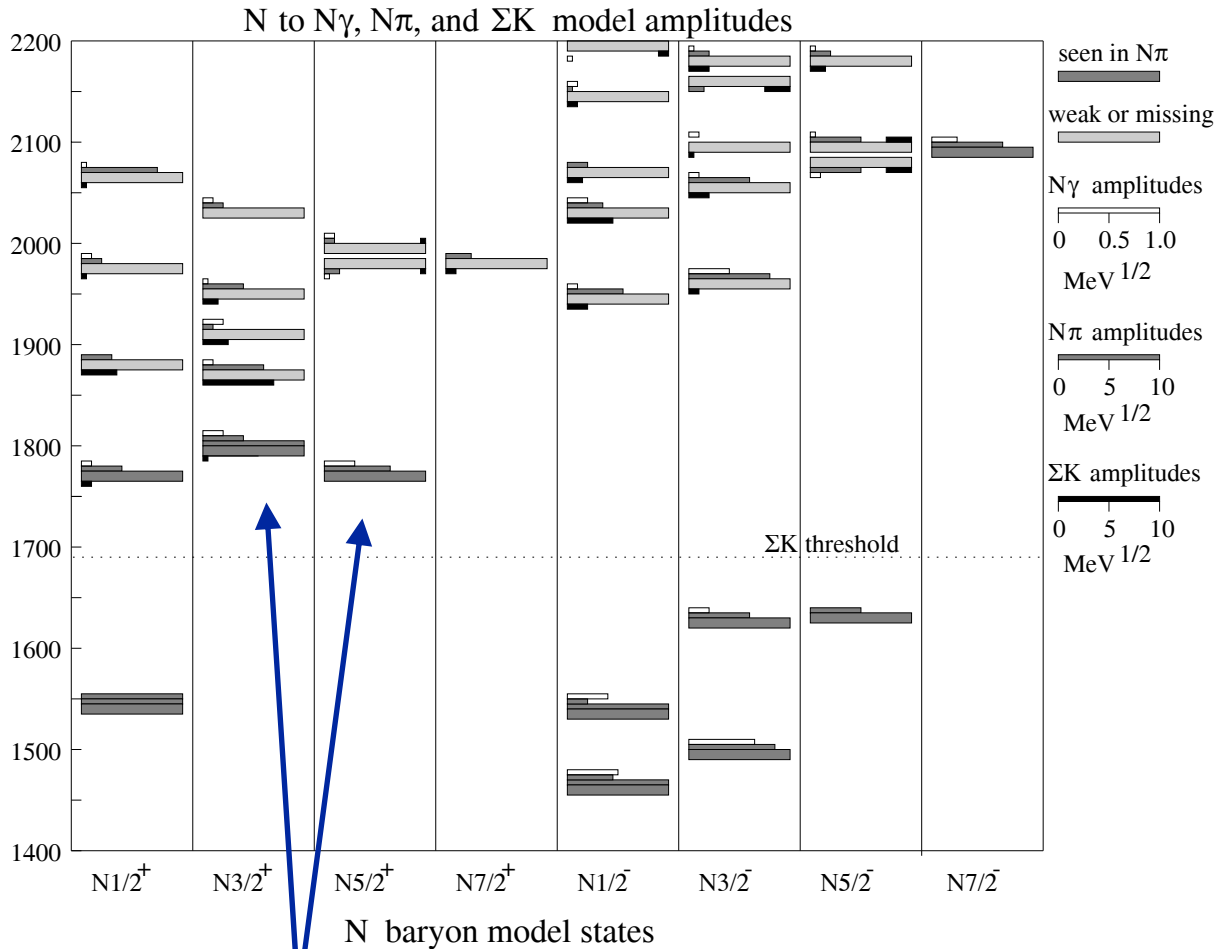
- *S. Basak, J. Dudek, R. Edwards, G. Fleming, U. Heller, J. Juge, A. Lichtl, N. Mathur, C. Morningstar, DGR, I. Sato, S. Wallace*



Spectroscopy and QCD

- Spectroscopy is classic tool for gleaning information about structure of theory
- Both experimental and ab initio N^* and Hybrid programs aim at discovering effective degrees of freedom of QCD, and resolving competing low-energy models:
 - Valence quark models with harmonic-oscillator potential cannot easily accommodate **Roper** and **$\Lambda(1405)$**
 - Quark models predict a richer spectrum of missing resonances than observed experimentally
 - **Diquark-quark picture of baryons** – *fewer degrees of freedom*
 - **Chiral soliton models**

Isospin-1/2 N* Sector



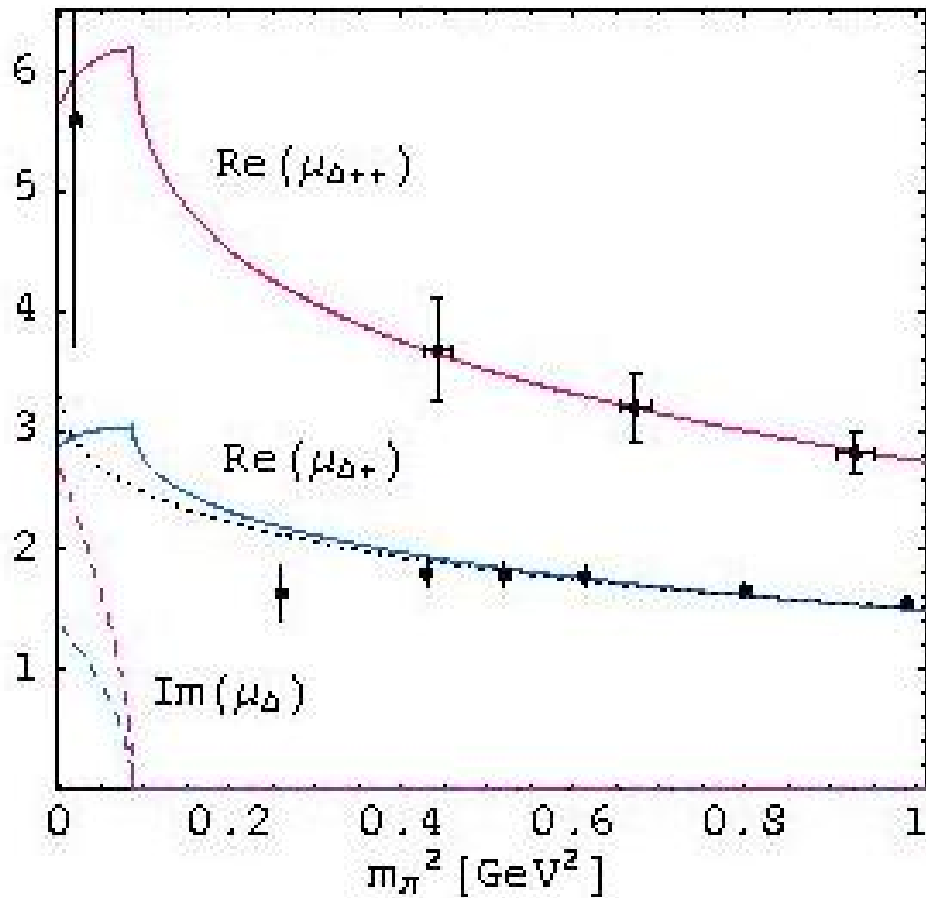
*Capstick and Roberts,
 PRD58 (1998) 074011*

Note ordering....



Spectroscopy review - I

One baryon resonance studied for many years.....

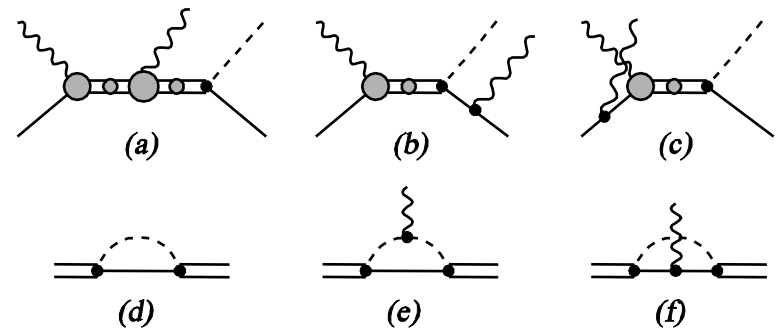


Lattice points from

Leinweber (1992)

Cloet, Leinweber, Thomas (2003)

Lee *et al.* (2004)



Chiral calculations

Pascalutsa, Vanderhaeghen (2004)

Thomas, Young (...)

Spectroscopy Review - II

- First generation calculations – largely for quarks masses around that of *strange quark*;
- Computations performed in quenched approximation to QCD.
- Masses of lowest lying baryon of each parity for $J = 1/2, 3/2$ for $N, \Delta, \Sigma, \Lambda$
- Computations performed using a variety of fermion actions – different $O(a)$ discretisation errors. But only limited investigation of systematic errors.
- Computations use simple three-quark interpolating fields - *insensitive to excited glue, multiquark states*



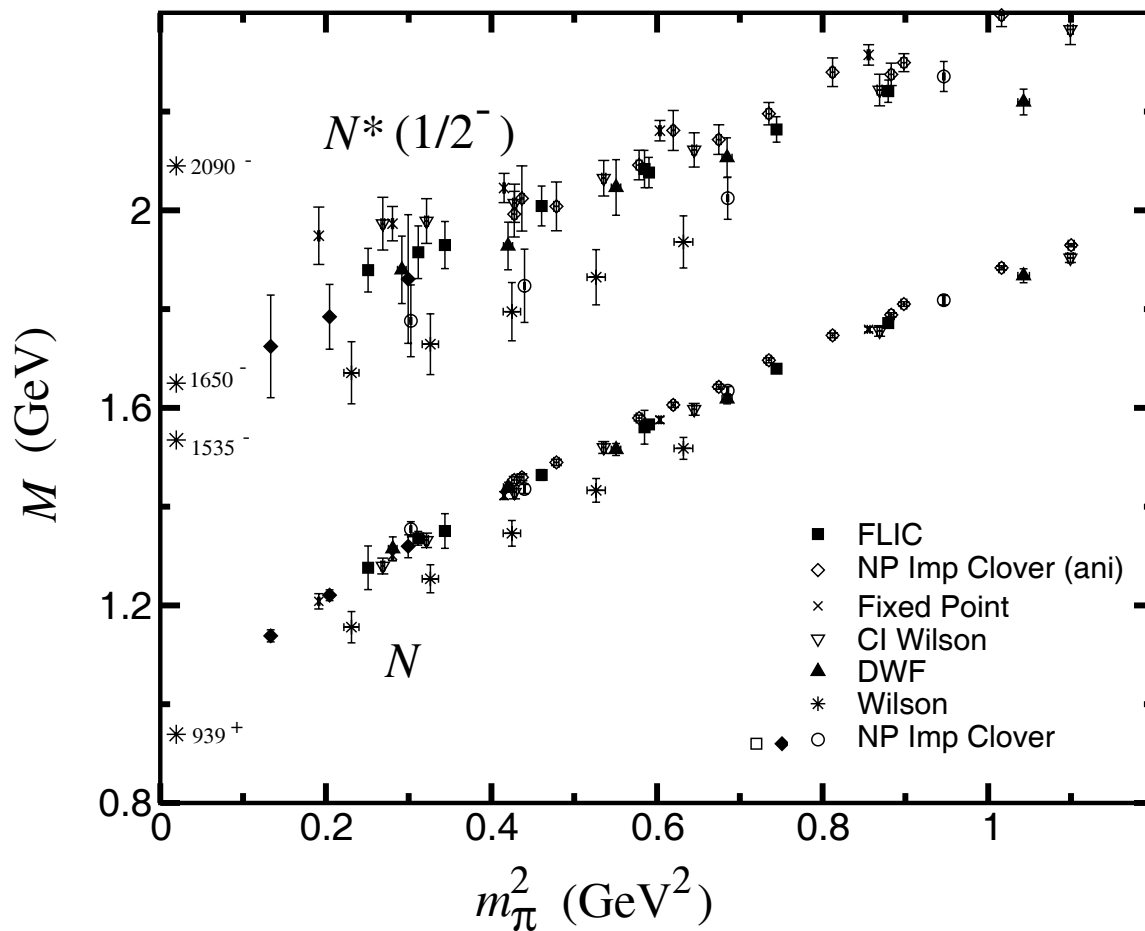
Parity Partner - I

Different symbols



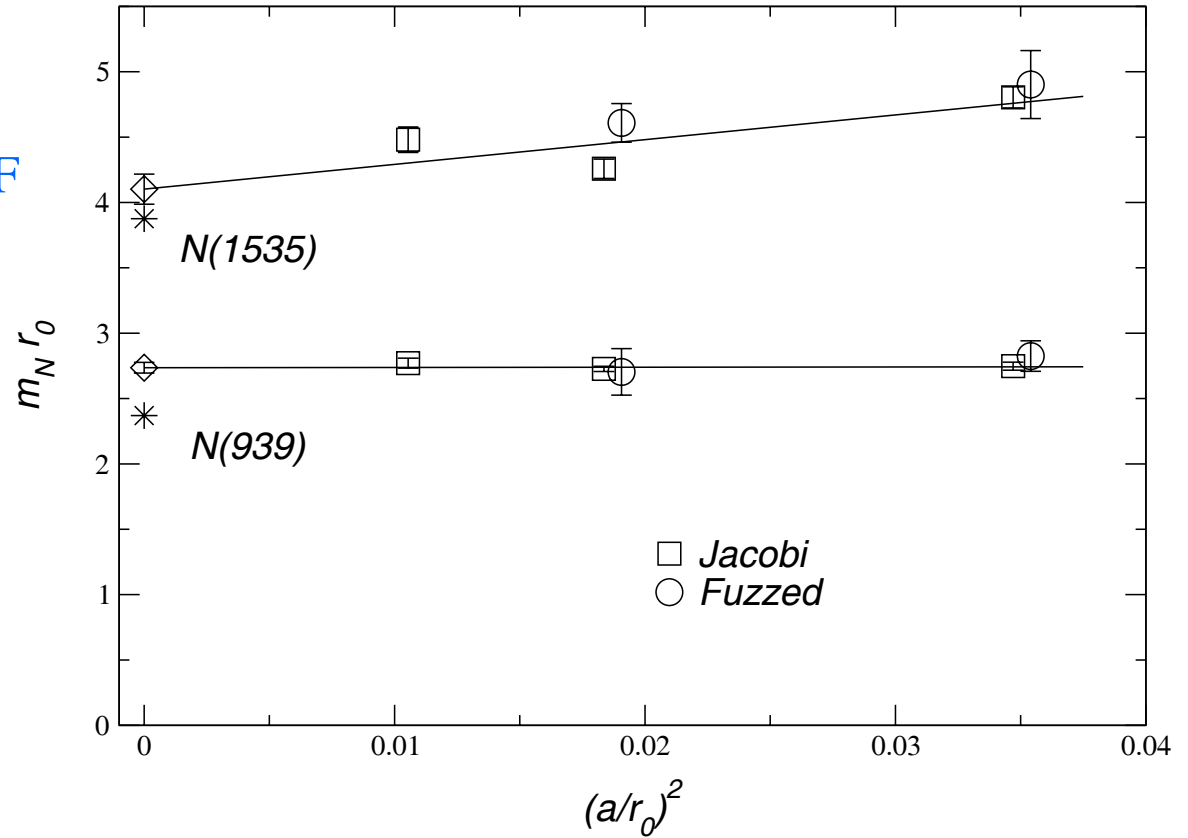
Different actions

Nucl-th/0406032



Parity partner - II

LHPC/UKQCD/QCDSF
PLB532, 63 (2002)

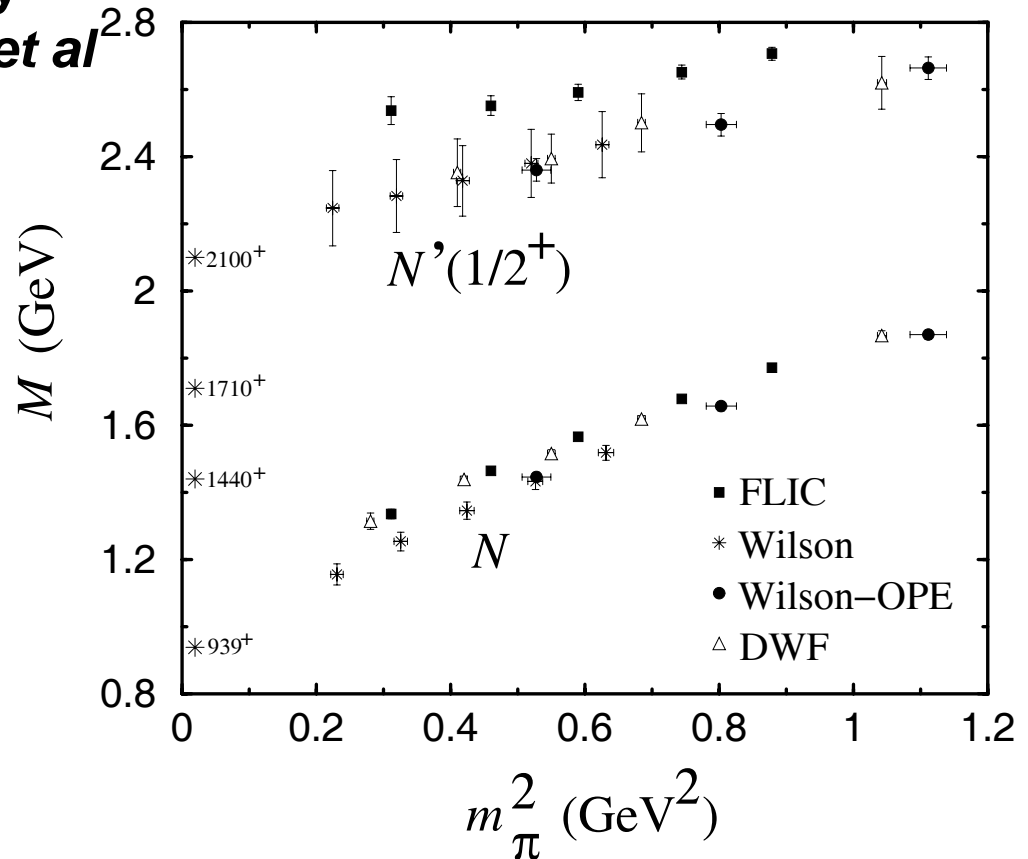


Nucleon Radial Excitation

Comparison of results by
Adelaide group, Zanotti *et al*
PRD67, 114506 (2003).

$$O_{N'} = (uCd)\gamma_5 u$$

“Bad” nucleon
operator – vanishes in
NR limit.

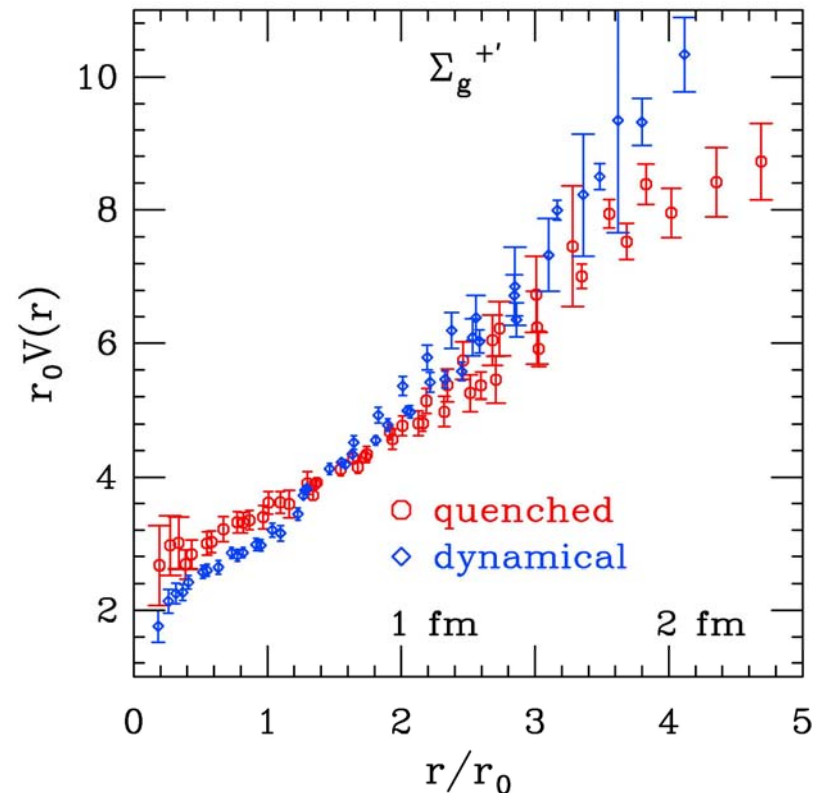


Need to extract radial and higher spin excitations

Hybrid Resonance Spectrum - I

- Spectrum of hybrids composed of heavy quarks has been important lattice focus
 - Adiabatic Potentials
 - Spectrum of charmonium
- 1 GeV Excitation Energy associated with excited string
 - 1⁺ lightest hybrid

MILC Collaboration

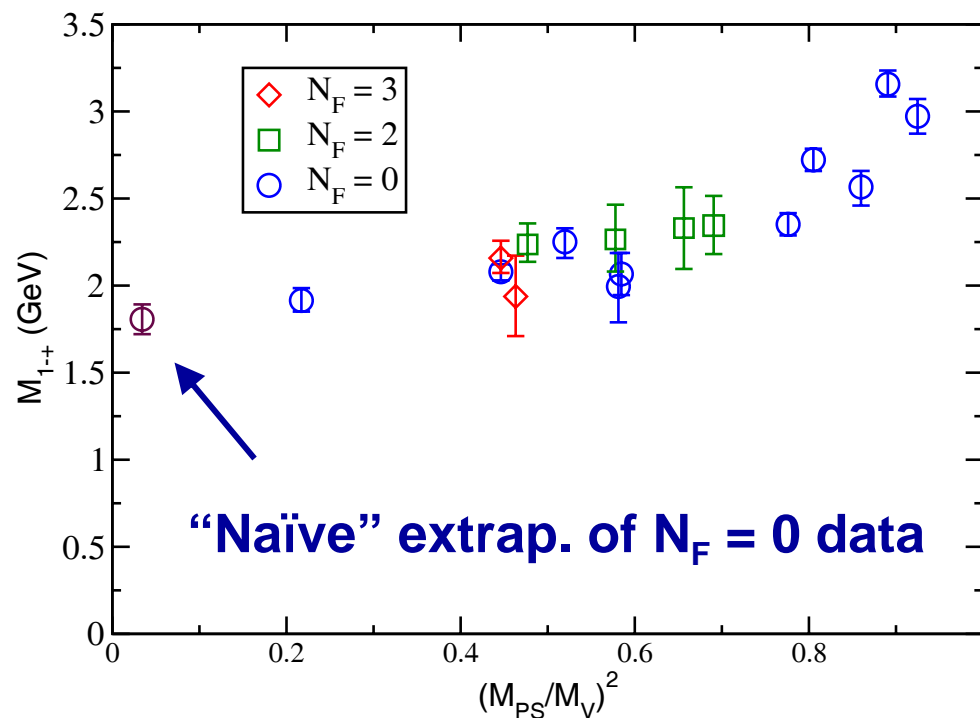


Hybrid Resonance Spectrum - II

- Does picture of hybrids obtained in charmonium sector persist to light-quark sector?

Compilation of data for mass of 1^{++} from *MILC Collaboration*

Scale set by string tension



Energies from correlation functions

- Stationary state energies can be extracted from asymptotic decay rate of temporal correlations of the fields (in the imaginary time formalism)
- Spectral representation of a simple correlation function
 - assume transfer matrix, ignore temporal boundary conditions
 - focus only on one time ordering

insert complete set of energy eigenstates

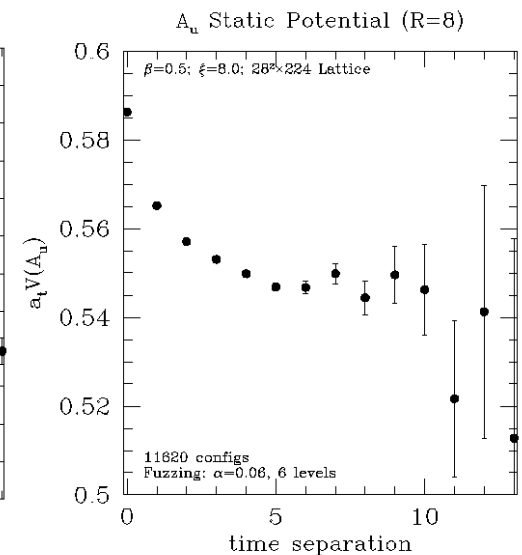
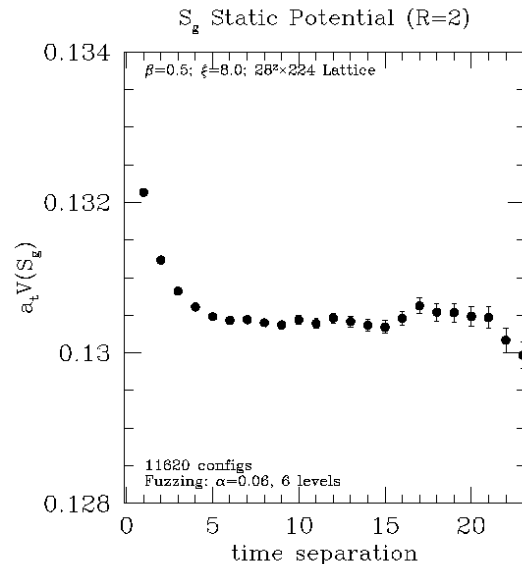
$$\begin{aligned} C(t) = \langle 0 | \phi(t)\phi(0) | 0 \rangle &= \sum_n \langle 0 | e^{Ht}\phi(0)e^{-Ht} | n \rangle \langle n | \phi(0) | 0 \rangle \\ &= |\langle n | \phi(0) | 0 \rangle|^2 e^{-E_n t} = \sum_n A_n e^{-E_n t} \end{aligned}$$

- *Extract lowest energy and amplitude as $t \rightarrow \infty$*

Effective mass

- The “effective mass” is given by $m_{\text{eff}}(t) = \ln \left(\frac{C(t)}{C(t+1)} \right)$
- The effective mass tends to the energy) asymptotically
- Effective mass plot is convenient visual tool to see signal extraction

- seen as a plateau
- Plateau sets in quickly for good operator
- Excited-state contamination before plateau



Principal correlators

- Extracting excited-state energies described in **C. Michael, NPB 259, 58 (1985)** and **Luscher and Wolff, NPB 339, 222 (1990)**
- Can be viewed as exploiting the *variational method*
- Given $N \times N$ correlator matrix $C_{\alpha\beta}(t) = \langle 0 | O_\alpha(t) O_\beta(0) | 0 \rangle$, one defines the N *principal correlators* $\lambda_i(t, t_0)$ as the eigenvalues of

$$C^{-1/2}(t_0)C(t)C^{-1/2}(t_0)$$

where t_0 (the time defining the “metric”) is small

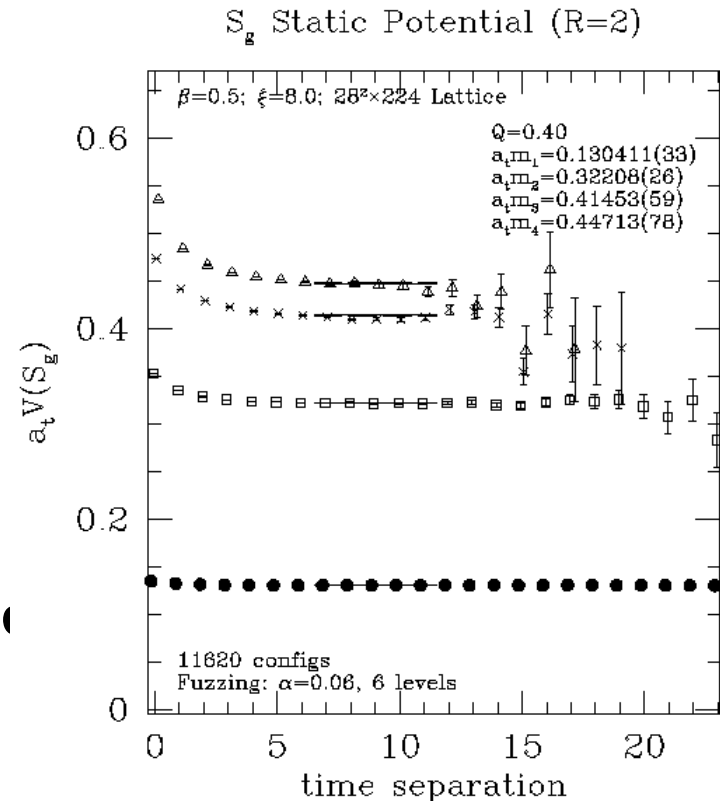
- Principal effective masses defined from correlators plateau to lowest-lying energies

$$\lambda_i(t, t_0) \rightarrow e^{-E_i(t-t_0)} \left(1 + O(e^{-\Delta E(t-t_0)}) \right)$$

$$\min (E_n - E_i)$$

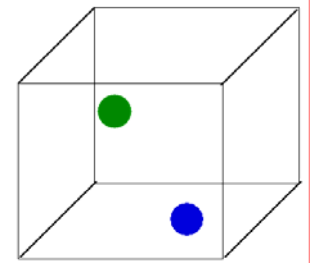

Principal effective masses

- Perform single-exponential fit to each principal correlator to extract spectrum! *Can use multi-exponentials to minimize sensitivity to excited states*
- Note that principal effective masses (as functions of time) can cross, approach asymptotic behaviour from below
- Final results are *independent of t_0* , but choosing larger values of this reference time can introduce larger errors

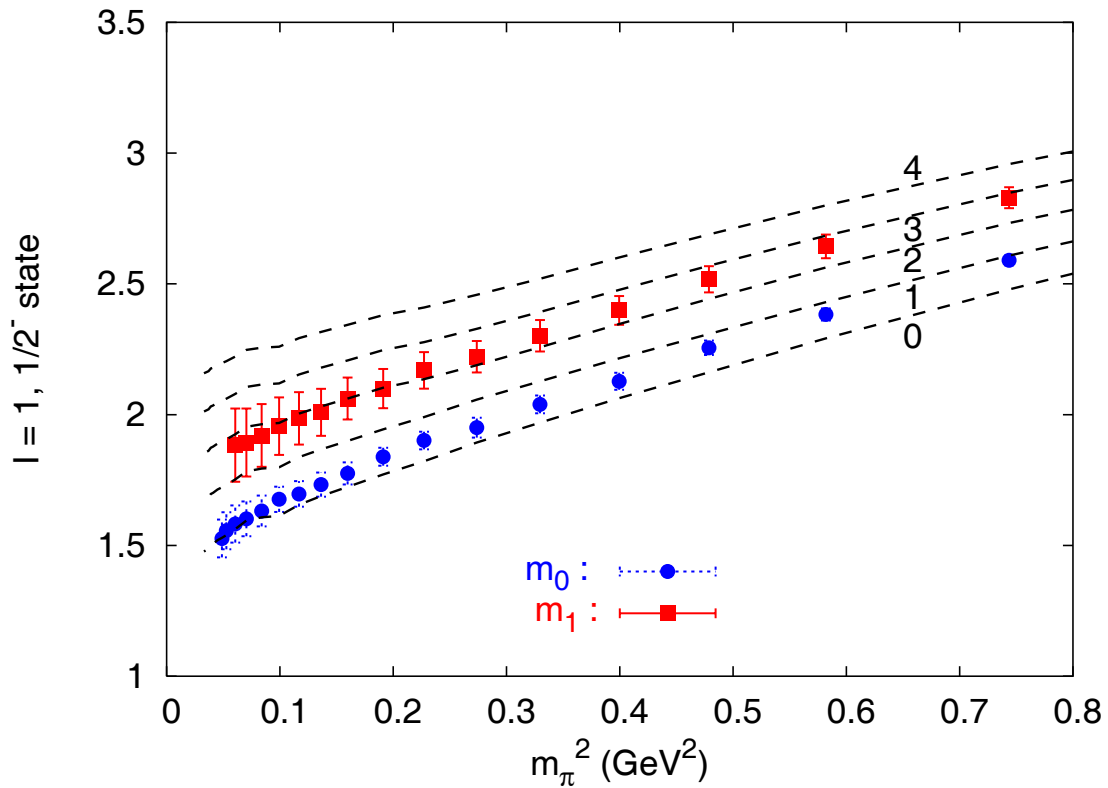


Unstable Particles

- computations done in a periodic box
 - momenta quantized
 - discrete energy spectrum of stationary states → single hadron, 2 hadron,...
- scattering phase shifts → resonance masses, widths (in principle) deduced from finite-box spectrum
 - B. DeWitt, PR **103**, 1565 (1956) (sphere)
 - M. Luscher, NPB**364**, 237 (1991) (cube)
- Two-particle states and resonances identified by examining behaviour of energies in finite volume
 - **Resonances with milder volume dependence**



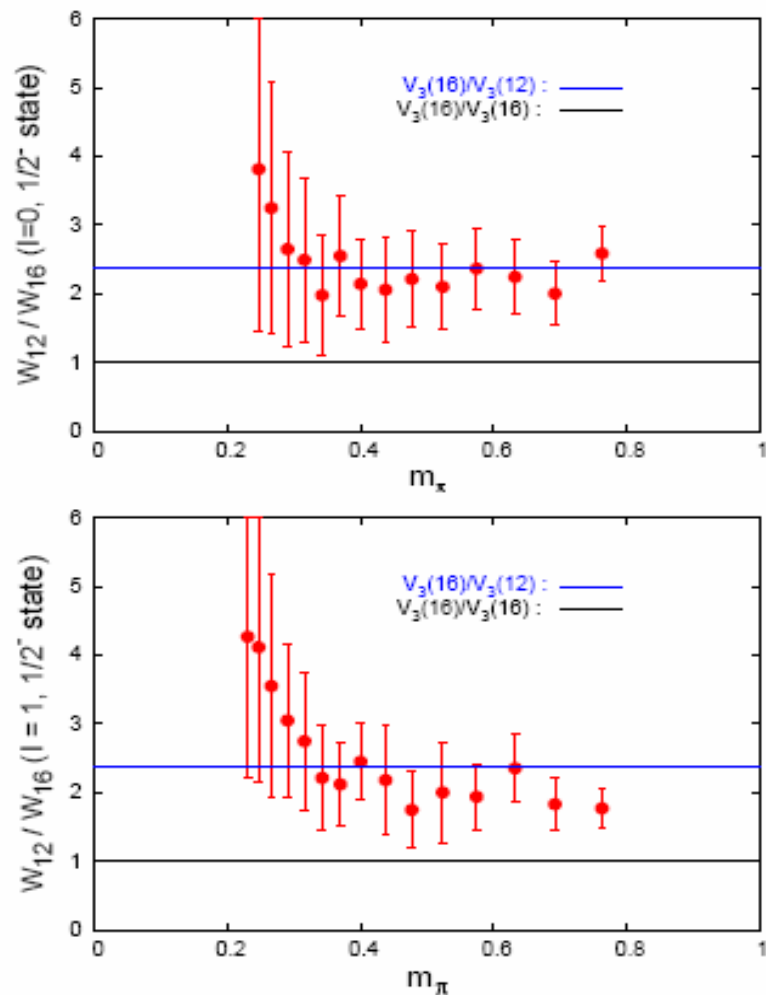
Resonance or scattering state - I



Mathur *et al*,
Phys.Rev. D70
(2004) 074508

Fit to ground and first-excited states of $I=1$ pentaquark correlator, together with scattering energies

Resonance or scattering state - II

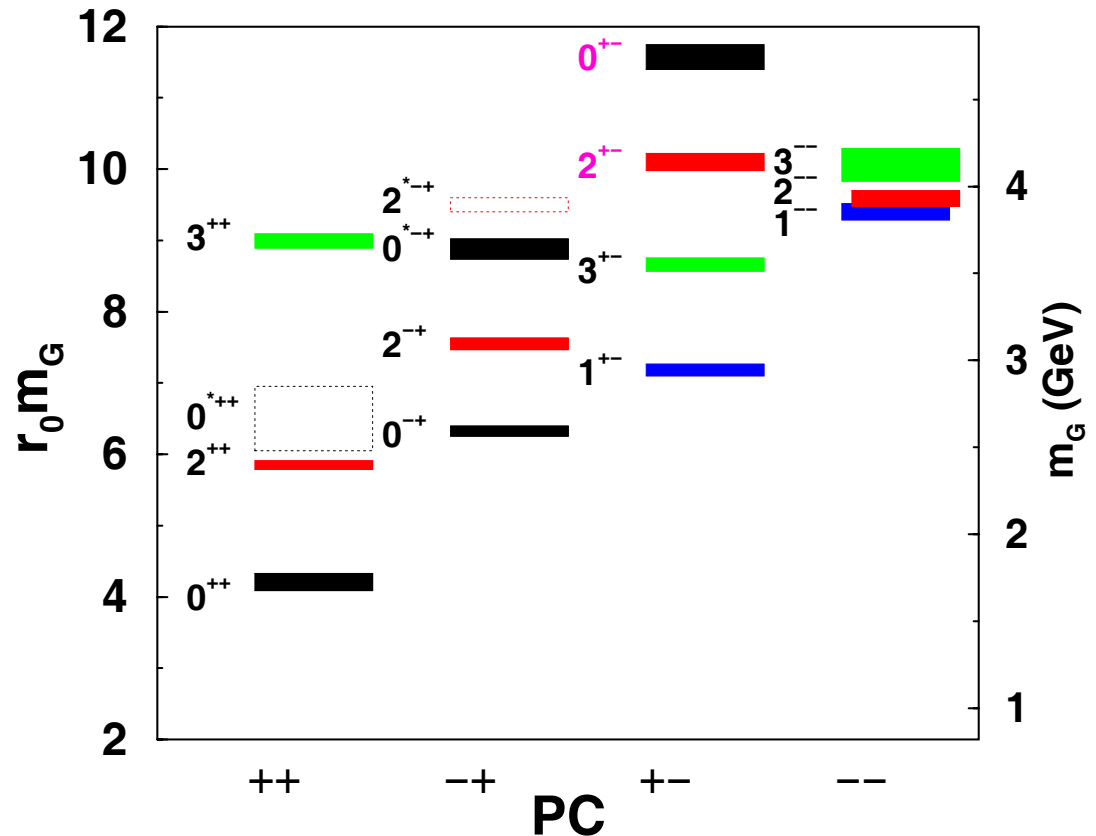


Two-particle amplitude expected to have different volume dependence to single-particle amplitude – *Amplitudes more consistent with two-particle state*

Variational methods: glueballs

Morningstar and Peardon
PRD60, 034509

- Quenched glueball calculations provide road-map
- Variational methods
- Large matrix of correlators
- Method more demanding for particles containing quarks



Baryons: Cubic Group of Lattice

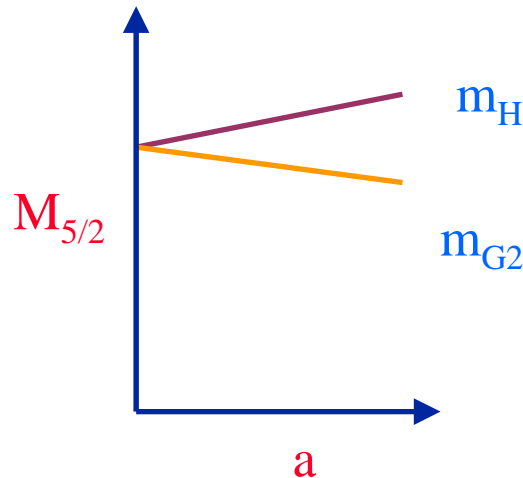
- Rotations restricted to isotropic cubic lattice form **Octohedral group O**
 - O has 24 elements
 - O has five conjugacy classes, hence five single-valued representations, **A₁(1), A₂(1), E(2), T₁(3), T₂(3)**
 - Under elements R of O, operator lying in Λ irreducible representation transforms as (λ is row)
$$U(R)O_{\lambda}^{(\Lambda)}U(R)^{\dagger} = \sum_{\lambda'} O_{\lambda'}^{(\Lambda)} D_{\lambda'\lambda}^{(\Lambda)}(R)$$
- **Spatial inversions** $I_s \rightarrow O_h$, with 48 elements; parities labelled g or u

Double Octohedral Group

- Has 48 elements
- Contains irreducible representations of O, together with **3 spinor irreps G_1, G_2, H** : R.C.Johnson, PLB114, 147 (82)

J	$n_{G_1}^J$	$n_{G_2}^J$	n_H^J
$\frac{1}{2}$	1	0	0
$\frac{3}{2}$	0	0	1
$\frac{5}{2}$	0	1	1
$\frac{7}{2}$	1	1	1
$\frac{9}{2}$	1	0	2
$\frac{11}{2}$	1	1	2
$\frac{13}{2}$	1	2	2
$\frac{15}{2}$	1	1	3
$\frac{17}{2}$	2	1	3

Note that states with $J \geq 5/2$ lie in representations with lower spins.



Spins identified from degeneracies in continuum limit

Baryon Fields and operators: design

- States are classified according to
 - Flavor structure, **F** *S. Basak et al., PRD72:074501,2005*
 - Parity *PRD72:094506,2005*
 - Total Spin, **J**, and Helicity, or third component of spin, **J_z**.
- In lattice computation, need to construct operators $B_i^{\Lambda\lambda F}(t, \mathbf{x})$ transforming irreducibly under **lattice symmetries** Λ, λ and form correlation matrix

$$C_{ij}^{\Lambda\lambda F}(t) = \sum_{\vec{x}} \langle 0 | B_i^{\Lambda\lambda F}(t, \vec{x}) \bar{B}_j^{\Lambda\lambda F}(0) | 0 \rangle$$

- Begin by considering flavor and orbital radial structure

Minimize Propagator Inversions

Flavour, Orbital and Radial Structure

- Classify states according to ***SU(2) Flavour***
- Build up ***radial structure*** using displacement operators **D**
- Provide our elemental **$B_i^F(t, \mathbf{x})$** operators having correct ***flavour and colour structure***

Quark-diquark

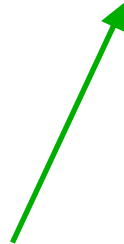


Illustration	Name	Explicit form ($ i \neq j \neq k $)
	single-site	$\phi_{ABC}^F \varepsilon_{abc} \tilde{\psi}_{A\alpha\alpha} \tilde{\psi}_{Bb\beta} \tilde{\psi}_{Cc\gamma}$
	singly-displaced	$\phi_{ABC}^F \varepsilon_{abc} \tilde{\psi}_{A\alpha\alpha} \tilde{\psi}_{Bb\beta} (\tilde{D}_j^{(p)} \tilde{\psi})_{Cc\gamma}$
	doubly-displaced-I	$\phi_{ABC}^F \varepsilon_{abc} \tilde{\psi}_{A\alpha\alpha} (\tilde{D}_{-j}^{(p)} \tilde{\psi})_{Bb\beta} (\tilde{D}_j^{(p)} \tilde{\psi})_{Cc\gamma}$
	doubly-displaced-L	$\phi_{ABC}^F \varepsilon_{abc} \tilde{\psi}_{A\alpha\alpha} (\tilde{D}_j^{(p)} \tilde{\psi})_{Bb\beta} (\tilde{D}_k^{(p)} \tilde{\psi})_{Cc\gamma}$
	triply-displaced-T	$\phi_{ABC}^F \varepsilon_{abc} (\tilde{D}_{-j}^{(p)} \tilde{\psi})_{A\alpha\alpha} (\tilde{D}_j^{(p)} \tilde{\psi})_{Bb\beta} (\tilde{D}_k^{(p)} \tilde{\psi})_{Cc\gamma}$
	triply-displaced-O	$\phi_{ABC}^F \varepsilon_{abc} (\tilde{D}_i^{(p)} \tilde{\psi})_{A\alpha\alpha} (\tilde{D}_j^{(p)} \tilde{\psi})_{Bb\beta} (\tilde{D}_k^{(p)} \tilde{\psi})_{Cc\gamma}$

Spatial structure of operators

Final step: projection formula

- **Central formula**

$$B_i^{\Lambda\lambda F}(t, \vec{x}) = \frac{d_\Lambda}{g_{O_h^D}} \sum_{R \in O_h^D} D_{\lambda\lambda}^{(\Lambda)*}(R) U_R B_i^F(t, \vec{x}) U_R^\dagger$$

- **Group theory performed in Maple/C++ – requires explicit Dirac basis.**

Irrep	Δ, Ω	N	Σ, Ξ	Λ
G_{1g}	221	443	664	656
G_{1u}	221	443	664	656
G_{2g}	188	376	564	556
G_{2u}	188	376	564	556
H_g	418	809	1227	1209
H_u	418	809	1227	1209

Total numbers of operators assuming two different displacement lengths

Several weeks on 16-node cluster

- **Include multi-hadron operators**



Simple example: local nucleon operators...

- Elemental operators:

$$\Phi_{\alpha\beta\gamma} = \epsilon^{abc} (u_{\alpha}^a u_{\beta}^b d_{\gamma}^c - u_{\gamma}^a u_{\beta}^b d_{\alpha}^c)$$

- Constraints:

$$\Phi_{\alpha\beta\gamma} + \Phi_{\gamma\beta\alpha} = 0$$

$$\Phi_{\alpha\beta\gamma} + \Phi_{\beta\gamma\alpha} + \Phi_{\gamma\alpha\beta} = 0$$

- 20 linearly-independent operators
- 3 embeddings of G_{1g} , G_{1u} , 1 embedding each of H_g , H_u ; **NO G_2**
- Three linearly independent nucleon operators, and a single $N^{*3/2}$ operator**

$(u C \gamma_5 d) u$, $(u C d) \gamma_5 u$, Rarita-Schwinger projection

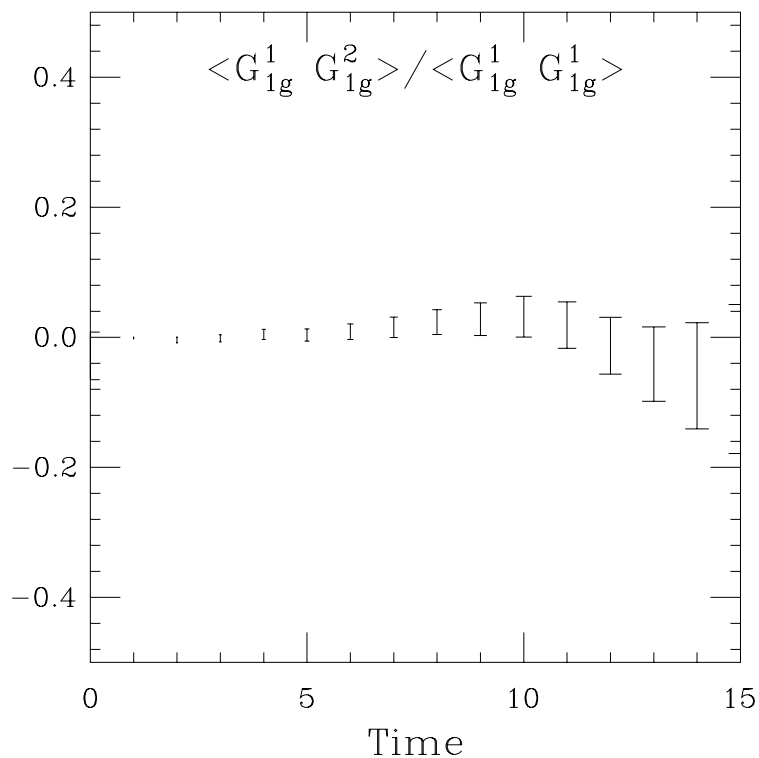
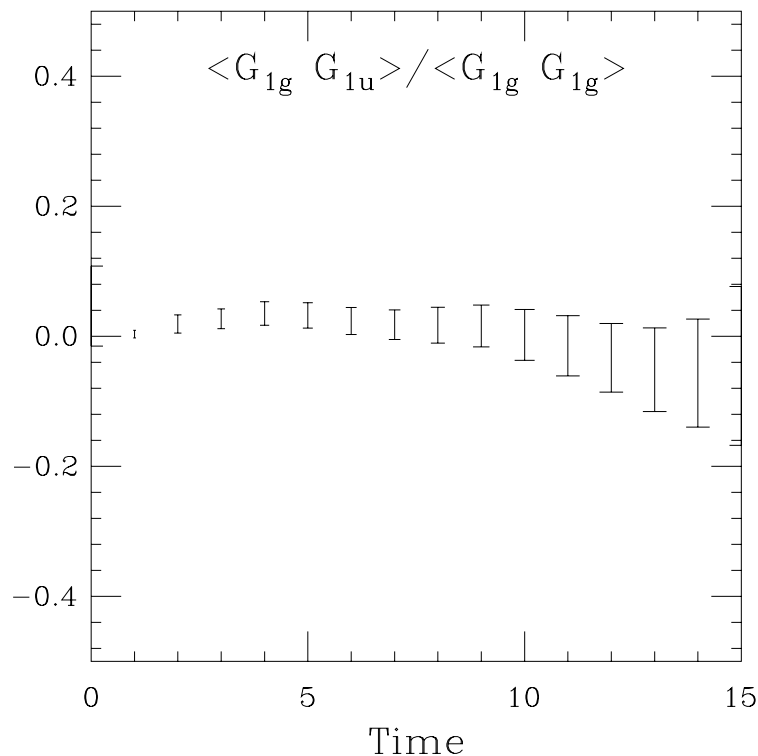
Implementation...

- On each configuration, form generalised baryon correlator

$$B_{\alpha\beta\gamma,\bar{\alpha}\bar{\beta}\bar{\gamma}} = \sum_{\vec{x}} \epsilon^{abc} \epsilon^{\bar{a}\bar{b}\bar{c}} A_{\alpha\bar{\alpha}}^{a\bar{a}}(x, 0) B_{\beta\bar{\beta}}^{b\bar{b}}(x, 0) C_{\gamma\bar{\gamma}}^{c\bar{c}}(x, 0)$$

- **Antisymmetrise** over α, γ - $I = 1/2$.
- **Symmetrise** over α and β – *Wick contractions*
- Form correlators according to table.
- Illustration: DWF valence propagators, on MILC Asqtad lattices:
 m_{π} ' 600 MeV
- **All-to-all propagators**

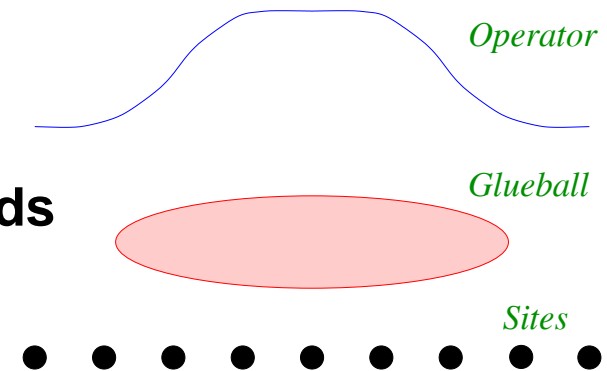
Orthogonality of States



- Different rows **orthogonal**
- Different irreps. **orthogonal**
- Different *embeddings* of same irrep **not orthogonal**

Smearing - I

- Statistical noise generally increases with temporal separation
- Effective masses associated with correlation functions of simple local fields do not reach a plateau before noise swamps the signal
 - need better operators
 - better operators have reduced couplings with higher-lying states
- Recipe for making better operators
 - Construct operators using *smear*ed fields
 - link variable smearing
 - quark field smearing



Implementation: Quark Smearing - I

- The quark smearing integral transform can be effected by the application of the smearing operator:

Adam Lichtl, Dublin '05

$$\tilde{\Psi}(x) = Q_{\sigma} \Psi(x)$$

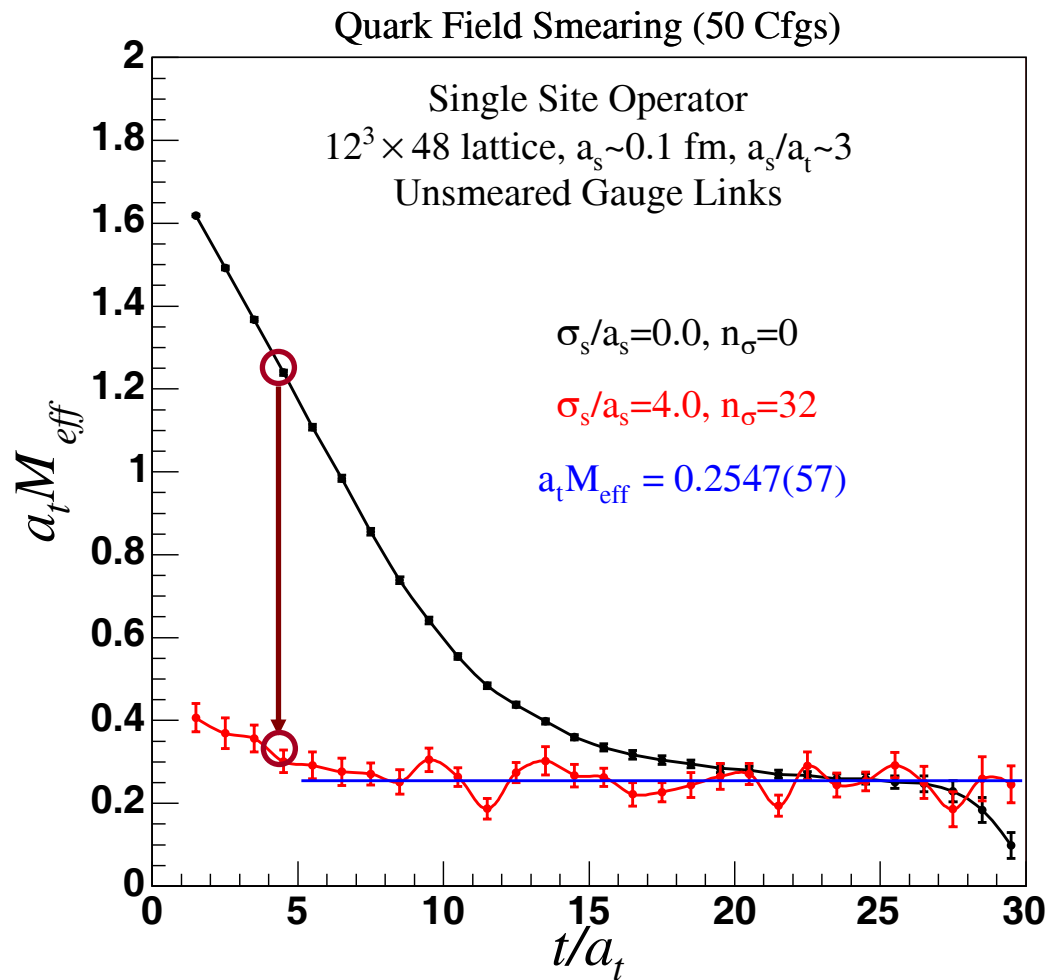
$$Q_{\sigma} = e^{\sigma^2 \nabla^2 / 2}$$

$$Q_{\sigma} = \lim_{n_{\sigma} \rightarrow \infty} (Q_{\sigma}^{n_{\sigma}}) = \lim_{n_{\sigma} \rightarrow \infty} \left(1 + \frac{\sigma^2 \Delta}{2n_{\sigma}}\right)^{n_{\sigma}}$$

- This attenuates the high momentum modes of the quark field → reduced contamination

The Chroma library uses: $Q(\sigma_s, n_{\sigma}) = \left(1 + \frac{\sigma_s^2 \Delta}{4n_{\sigma}}\right)^{n_{\sigma}}$

Implementation: Quark Smearing - II



Implementation: Link Smearing - I

- We see that link smearing plays a crucial role when we construct displaced operators
- **Stout link smearing** combines neighboring links in an analytic manner
 - Morningstar and Peardon, hep-lat/0311018

$$U_{\mu}^{(n+1)}(x) = e^{i\rho\Theta_{\mu}^{(n)}(x)}U_{\mu}^{(n)}(x)$$

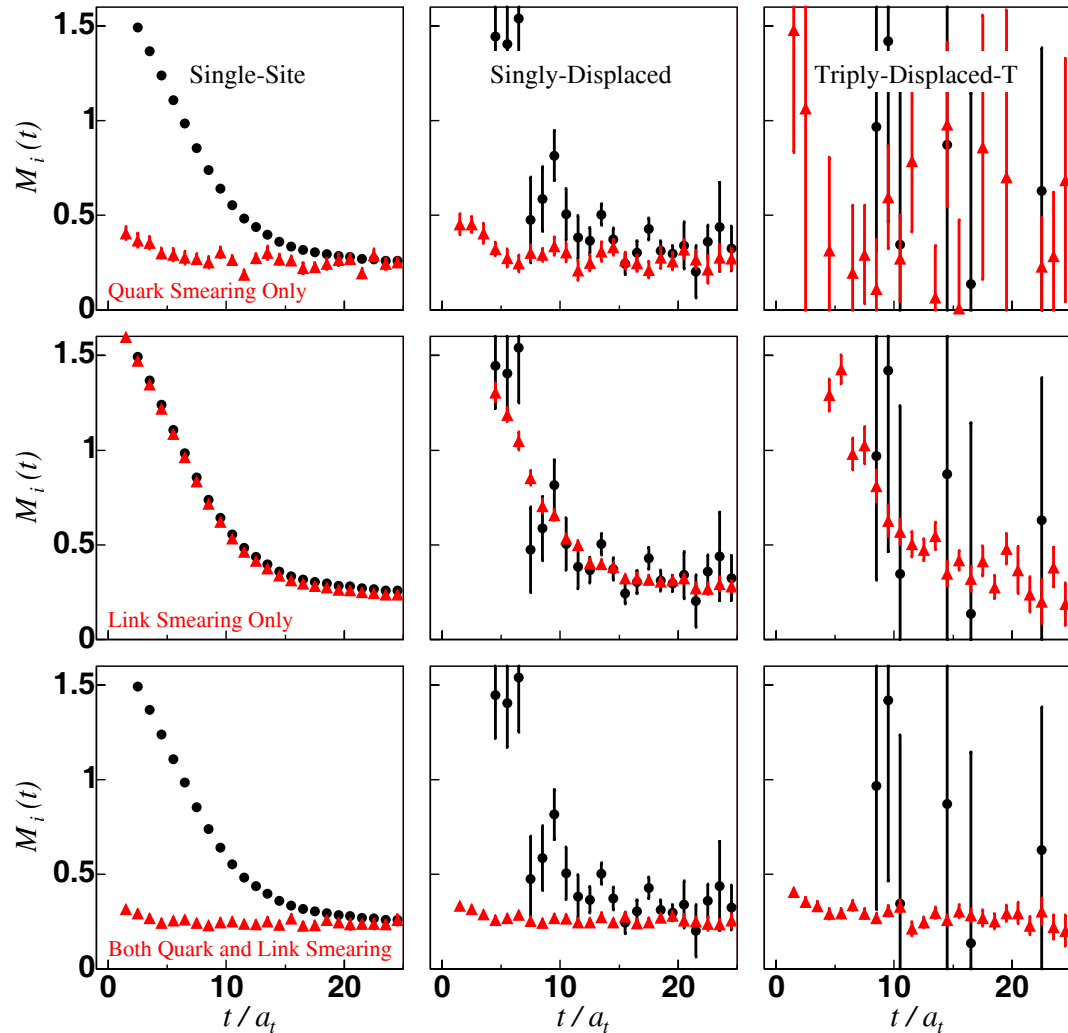
$$\Theta_{\mu}^{(n)} = \frac{i}{2}(\Omega_{\mu}^{\dagger}(x) - \Omega_{\mu}(x)) - \frac{i}{2N}\text{Tr}(\Omega_{\mu}^{\dagger}(x) - \Omega_{\mu}(x))$$

$$\Omega_{\mu}(x) = C_{\mu}(x)U_{\mu}^{\dagger}(x)$$

$$C_{\mu}(x) = \text{Sum of spatial staples from } x \text{ to } x + \hat{\mu}$$

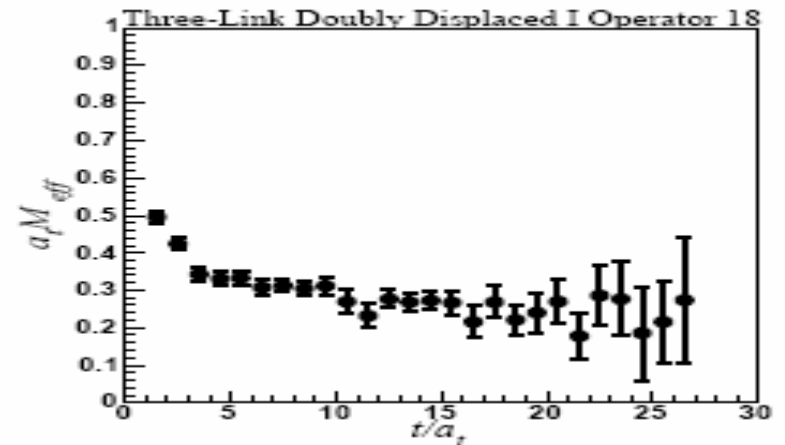
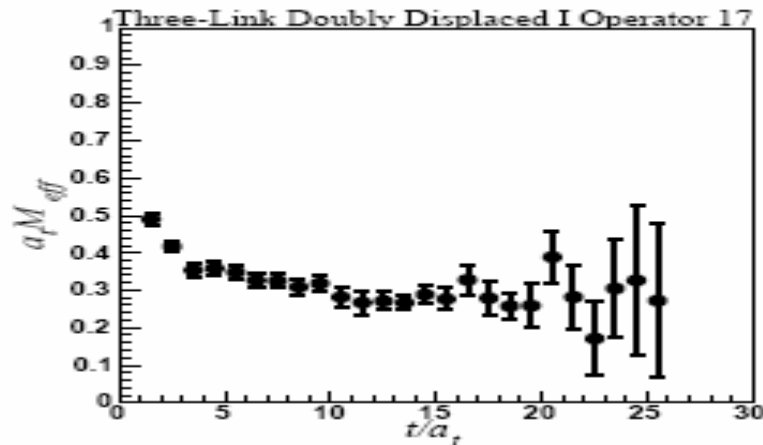
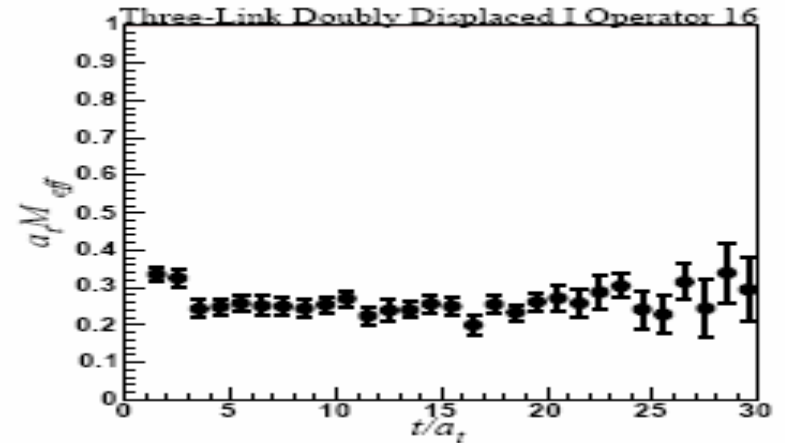
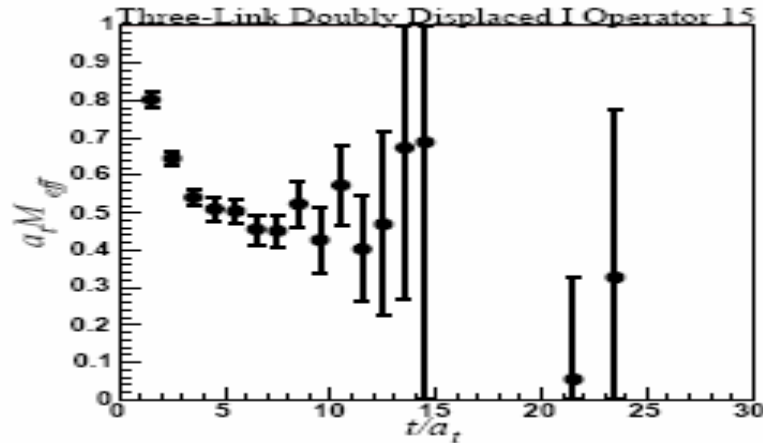
Implementation: Link Smearing - II

- **Quark field smearing**: reduces coupling to high-lying states.
- **Link smearing**: reduces variance in correlator.



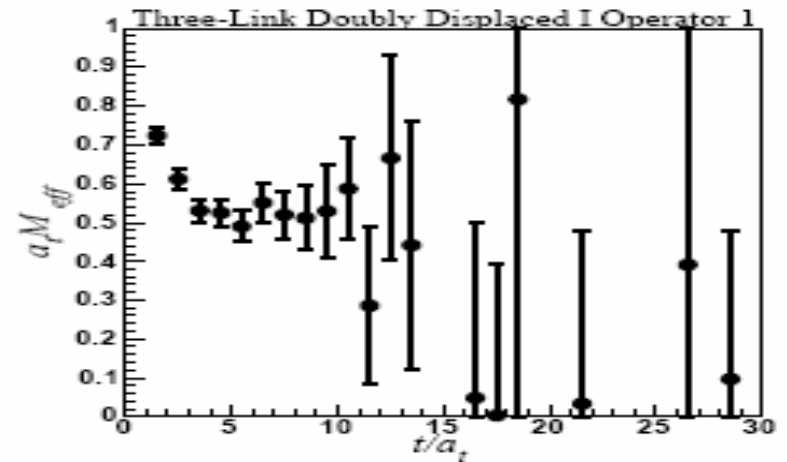
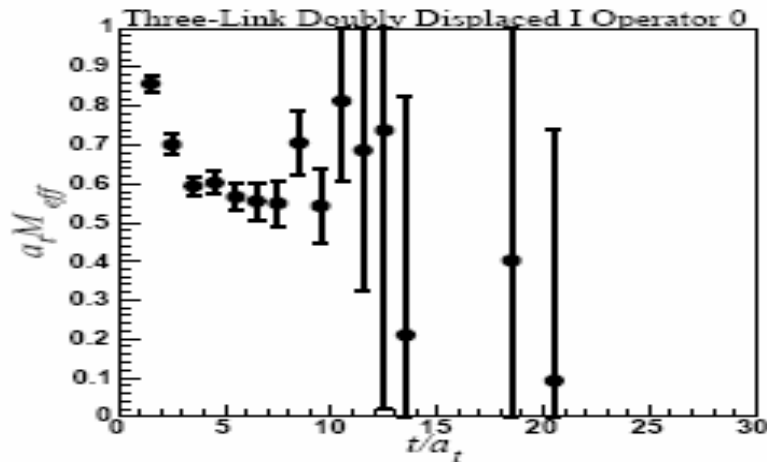
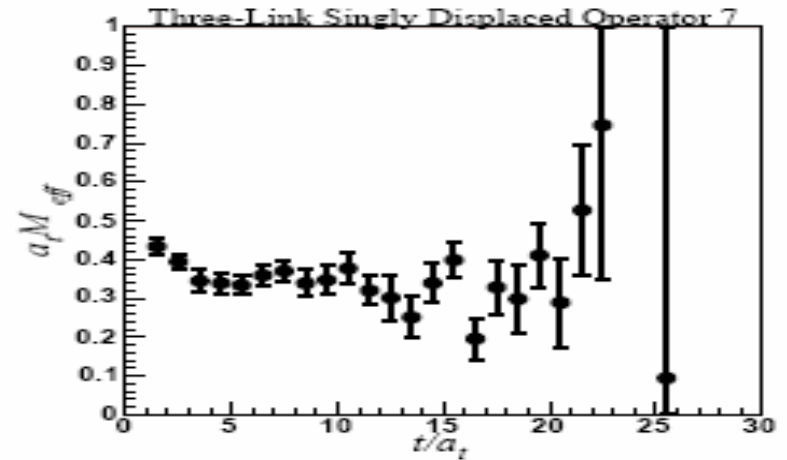
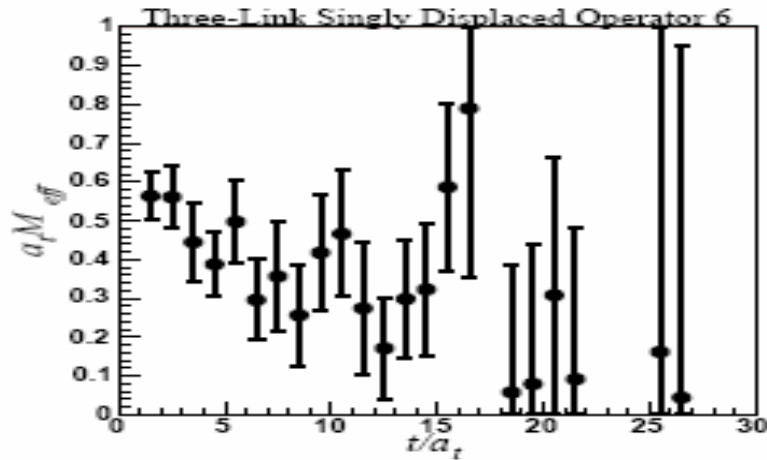
Implementation: Pruning - I

G_{1g}



Implementation - II

G_{2u}



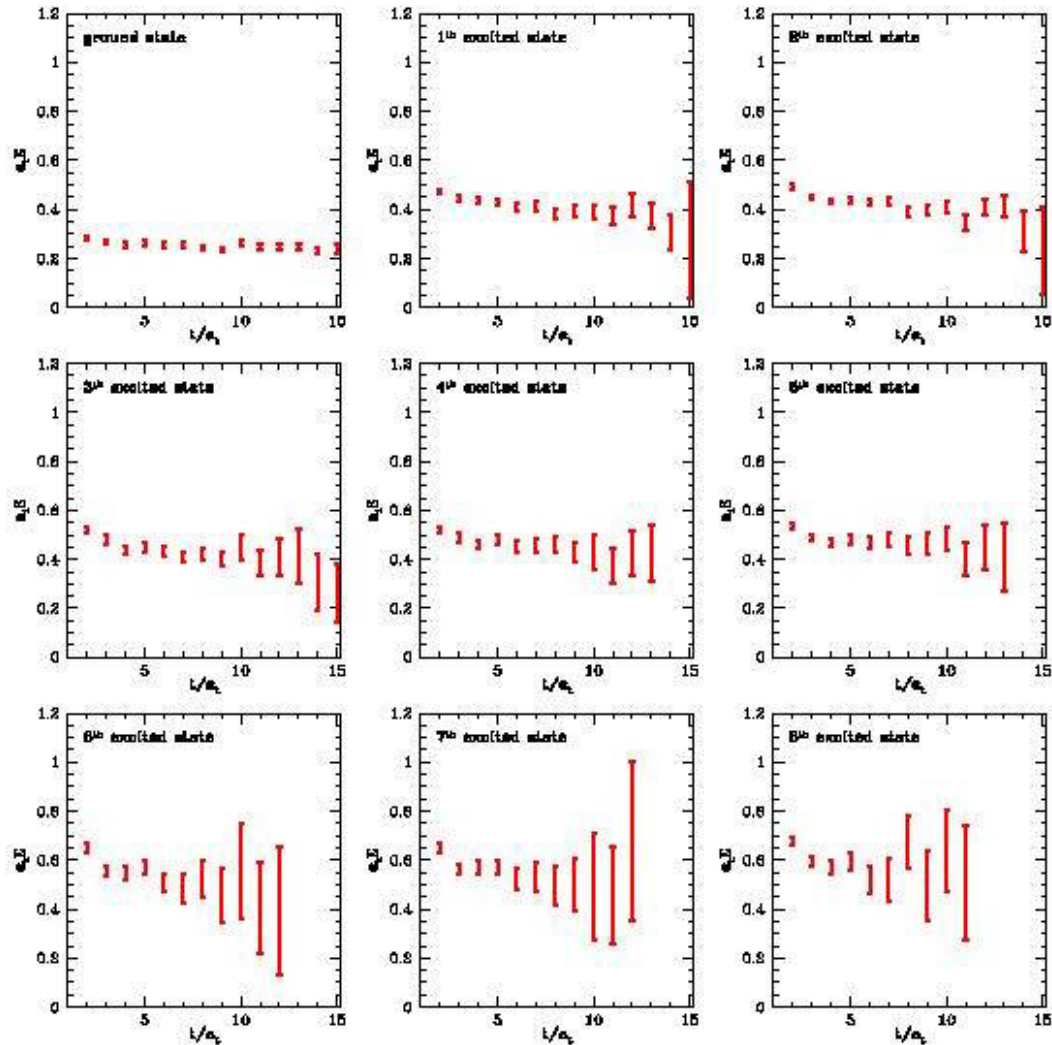
***A priori*, we do not know the “good” interpolating fields**



Principle Eigenvalues

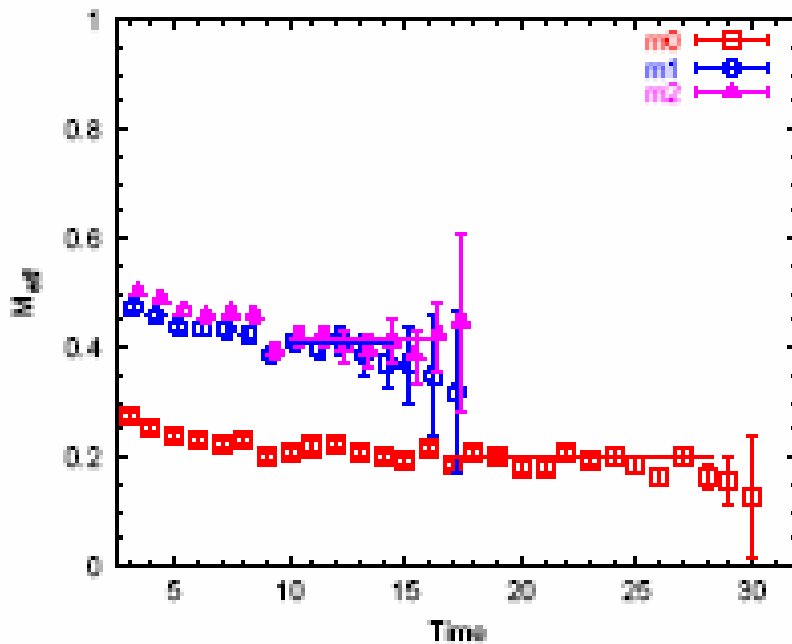
Extract up to 8
or 9 principle
correlators

cf **BGR**
Collaboration:
Gaussians of
different widths



Principle eigenvalues

- principal effective masses for small set of 10 operators



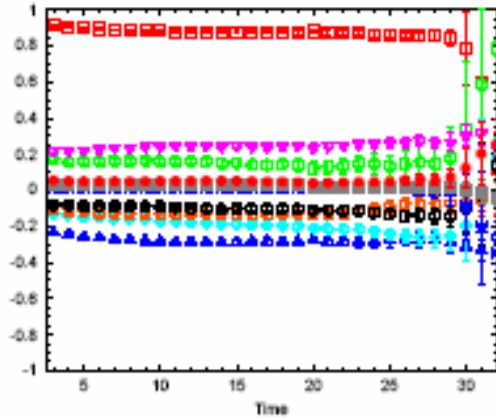
P-wave sources

Ikuro Sato, PhD, UMD (2005)

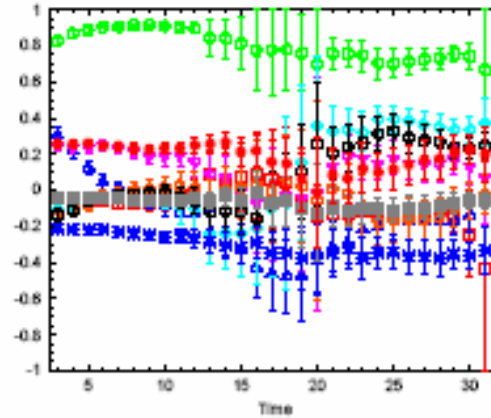
$16^3 \times 64$ lattice, anisotropic
Wilson with $\xi = 3.0$; $m_\pi =$
500 MeV

Eigenvectors....

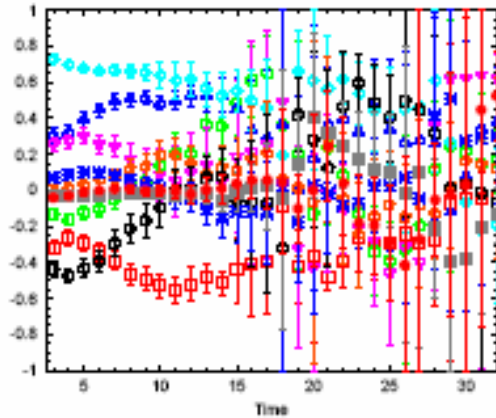
- Eigenvectors provide information regarding basis functions



(a) Ground state



(b) first-excited state

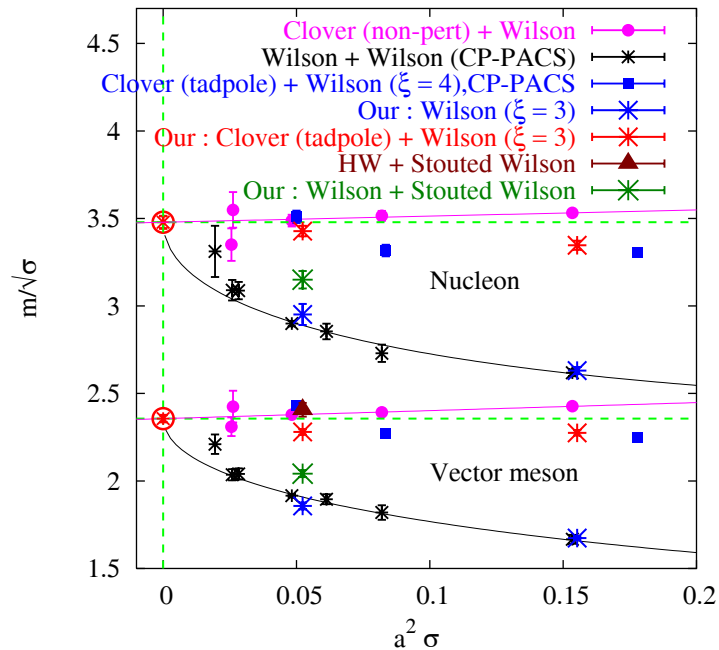


- G1g MA 1
- G1g MA 2
- G1g MAxMA T1 g1
- G1g MSxS T1 h1
- G1g MAxMA T1 h1
- G1g MAxMA T1 g3
- G1g MAxMA T1 g2
- G1g MSxS T1 g1
- G1g MSxS T1 h2

P-wave sources

Future Plans

- Generate dynamical anisotropic lattices
- Two lattice spacings – *spin identification*
- Three volumes – *Isolate two-particle contributions*
- Choice of action:
 - Require well-defined transfer matrix
 - ~~DWF or overlap~~
 - **Anisotropic clover action**



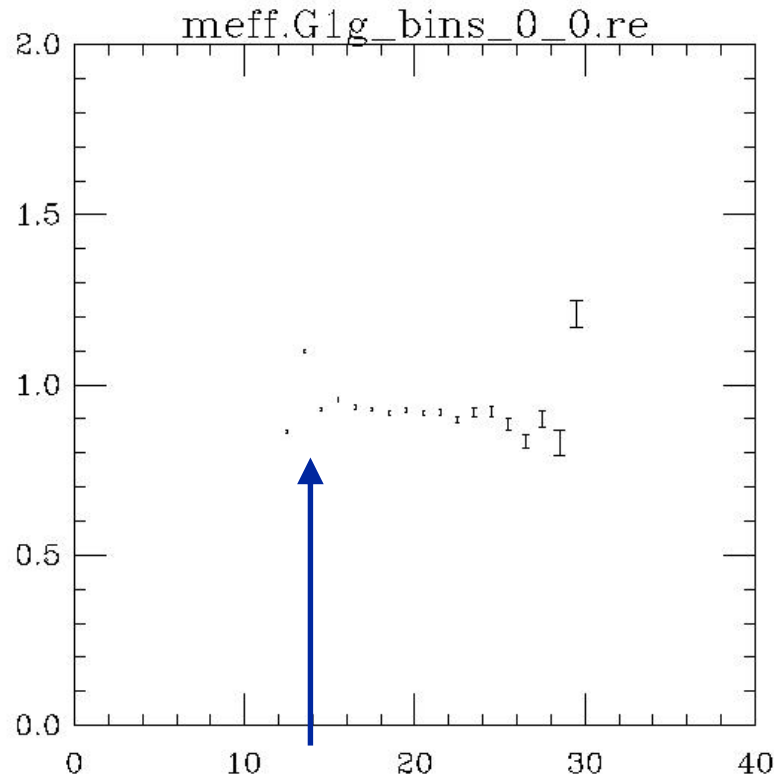
Choice of Action

- LHPC hadron structure studies employ DWF fermions on asqtad gauge background.
- DWF is $O(a)$ -improved – *crucial for study of matrix elements.*
- DWF is non-local in time and obscures interpretation of transfer matrix – *extremely problematic for resonance spectroscopy.*
- *Study of baryon resonance spectrum using all possible local operators*

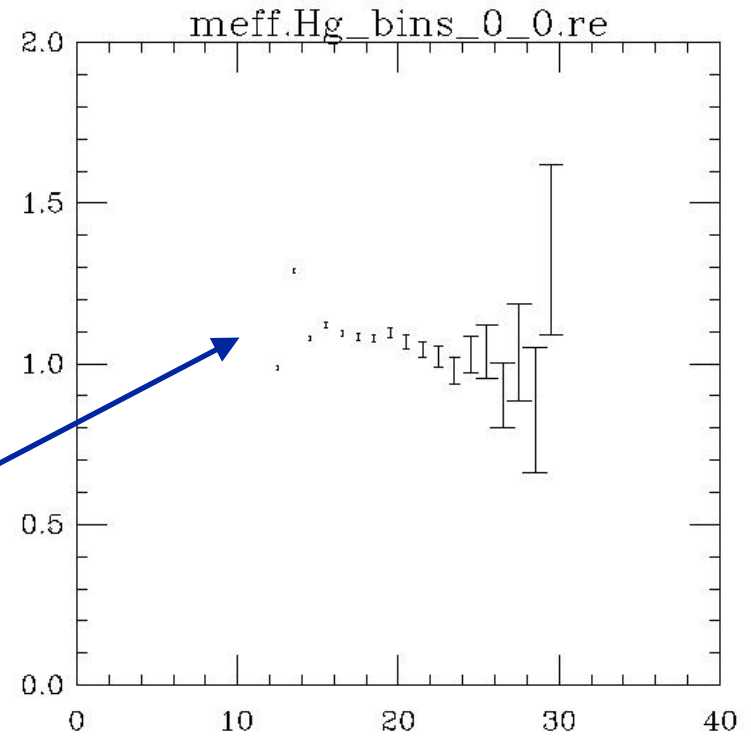


DWF - I

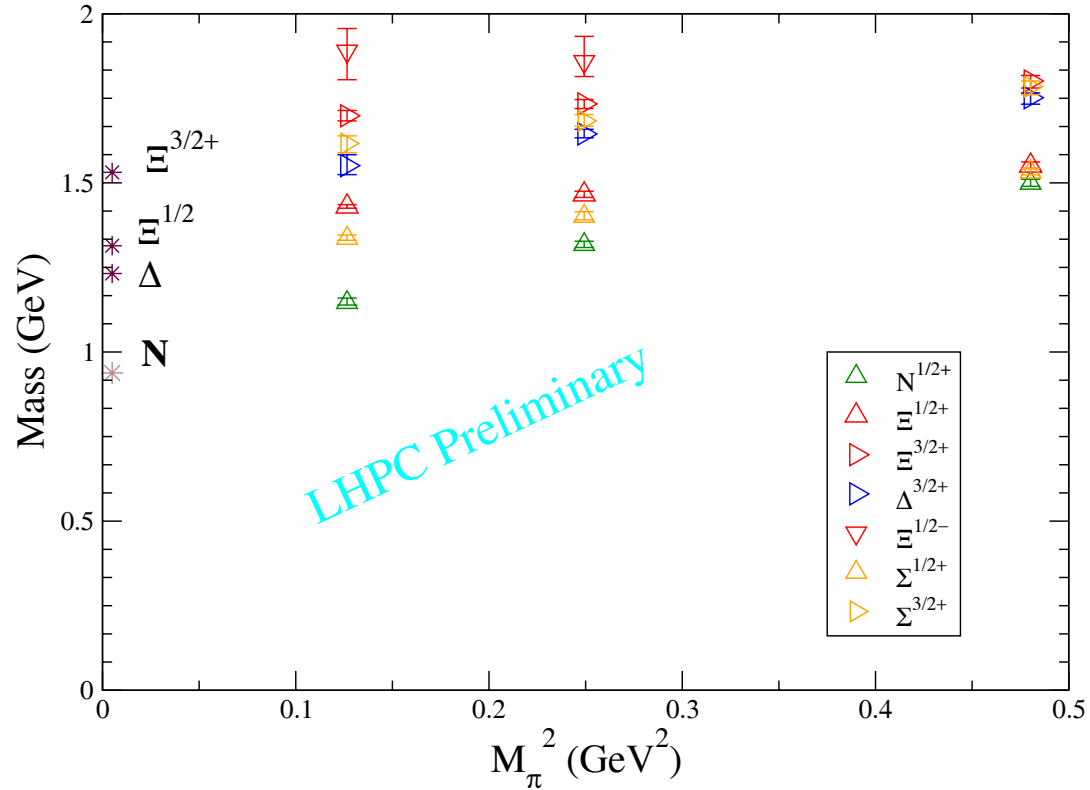
NPLQCD + LHPC Propagatos



Non-locality at short distance



DWF- II



Focus on cascades: *narrow width*

Conclusions and Perspective

- Exploration of resonance spectrum requires use of variational methods and development of suitable basis of operators
- Developed baryon operators lying in IR's of cubic group of the lattice – *essential to identify spin and parity*
- Operator construction is being extended to include **excited glue**, and to **hybrids**
- Proposal to generate dynamical, anisotropic clover lattices
- Aiming to establish efficacy of all-to-all methods – do not require choice of operator *Juge*
- For “well-established” resonances, can straightforwardly study radiative transitions –
 - $\gamma N \rightarrow \Delta$,
 - Radiative transitions in charmonium – *Dudek, Edwards, Mathur, DGR*