## Parallel Numerical Solutions of the QCD Bethe-Salpeter Equation II

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CMPT 851 Final Presentation

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- The purpose of the project is to model 2-body bound states in quantum chromodynamics (QCD).
- The wavefunction of a bound state can be obtained by solving the QCD Bethe-Salpeter equation:

$$\Gamma(p; P) = \int \frac{d^4q}{(2\pi)^4} K(q, p; P) S(q^+) \Gamma(q; P) S(q^-).$$

- This is an integral equation for the Bethe-Salpeter amplitude Γ(p; P), which is a 4 × 4 matrix.
- The equation can be made more numerically accessible by expanding the solution in various degrees of freedom.

#### Expansions

The first step is to expand the solution in Lorentz components (reducing the problem from an IE for a 4 × 4 solution to 4 coupled IEs for 1 × 1 solutions):

$$\Gamma(p; P) = \sum_{i=1}^{N} F_i(p; P) \widehat{\tau}_i(p; P).$$

 Then a Chebyshev decomposition can be performed to reduce 4D integration to a set of coupled 1D integrations:

$$F_i(p,P) = \sum_{j=0}^M U_j(u) \tilde{F}_{ij}(p^2,P^2).$$

 The remaining integration can be discretized, and an artifical eigenvalue can be inserted to allow the external momentum scale P to be fixed:

$$\lambda\left(P^{2}\right)\tilde{F}_{ij}\left(p^{2}\right)=\mathscr{K}\left(q^{2},p^{2},P^{2}\right)\cdot\tilde{F}_{ij}\left(q^{2}\right).$$

# Evaluation of $\mathscr{K}(P^2)$

- Evaluating the elements of  $\mathscr{K}(P^2)$  requires double integrations over the angular variables u. These integrations were attempted using both an adaptive Simpson quadrature rule and a Gauss-Legendre nonadaptive quadrature rule.
- Timings for adaptive Simpson quadrature (in seconds):



## Evaluation of $\mathscr{K}(P^2)$

■ Timings for Gauss-Legendre quadrature (in seconds):



The two methods produce results that are the same to within 1%:  $\max \left[ \frac{\mathscr{K}_{Simp} - \mathscr{K}_{GL}}{\mathscr{K}_{Simp}} \right] = 0.0033.$ 

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#### Parallelization of Problem

■ The problem can be visualized as evaluation of a 3D array:



■ Several methods were implemented on Socrates for the basic problem (32 grid size, both expansions truncated after one entry: N = M = 1). The following tables show minimum timings over 5 runs for 80 different values of P.

#### Parallelization of Problem

These two tables are for versions which parallelize the loop over P (Table A), and the loop over the elements of *K* (Table B).

A)	р	Time (seconds)	Speedup	] B)	р	Time (seconds)	Speedup
	1	161.522337	1		1	166.076348	1
	2	85.508760	1.8890		2	89.910633	1.8471
	4	49.311946	3.2755		4	53.421731	3.1088
	8	36.629304	4.4096		8	41.934506	3.9604

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### Parallelization of Problem

- The left table (C) makes both the loops over P and over the elements of *K* parfor loops.
- The right hand table (D) involves a single parallel loop over the elements of the full 3D array.

C)	р	Time (seconds)	Speedup	] D)	р	Time (seconds)	Speedup
	1	162.320255	1		1	163.824436	1
	2	85.886755	1.8899		2	87.518920	1.8719
	4	49.494977	3.2795		4	49.092138	3.3371
	8	36.357750	4.4645		8	34.331397	4.7719

#### Discussion

- For the basic problem, all these methods perform similarly.
- It is hard to draw any conclusions with only 8 workers available (maximum on Socrates).
- Note that versions A and C are functionally the same, given how workers are assigned.
- Version B likely scales slightly worse than the others due to overhead associated with setting up the loops.
- It is likely that version D will scale better than the others.
- Version D requires that the full problem be stored in memory (as a 3D array) to obtain the eigenvalues, and this may require too much memory with the full problem.

### Future Work

- Obviously the main extension is to use the full kernel *K* and solve the complete problem.
- Tests with the basic problem need to be done on a larger system to see which method actually scales the best.
- The convergence rates of the Chebyshev expansions also need to be examined.

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And finally, extension to 3-body problems.