

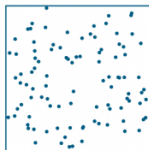
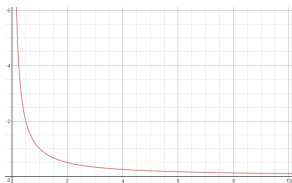
# A High-Order Fast Algorithm Approach for Computing Layer and Volume Potentials

J. Bevan, UIUC

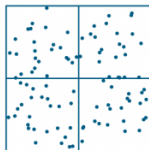
*CS 591 Seminar*  
*April, 2017*

# Motivation

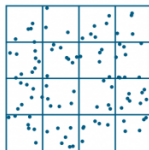
- ▶ Evaluation of potentials (derived velocity field, electric potential, etc.) is an important physical consideration
- ▶ Practical computational im[plementation] faces two challenges:  $\mathcal{O}(n^2)$  cost and integrable singularities



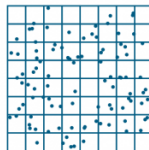
(a)  $d = 0$



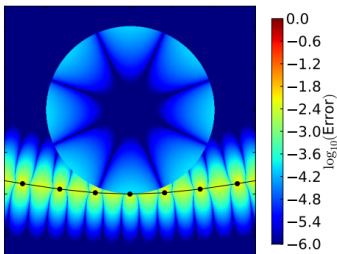
(b)  $d = 1$



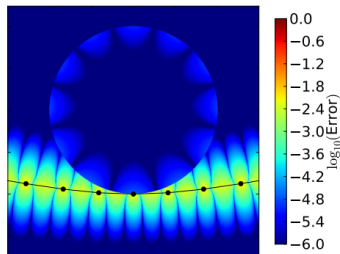
(c)  $d = 2$



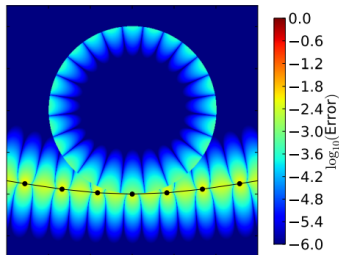
(d)  $d = 3$



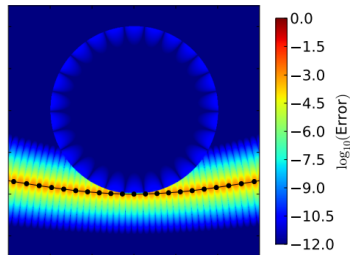
(a)  $p = 3$ ,  $N = 80$  quadrature nodes



(b)  $p = 6$ ,  $N = 80$  quadrature nodes



(c)  $p = 12$ ,  $N = 80$  quadrature nodes



(d)  $p = 12$ ,  $N = 240$  quadrature nodes

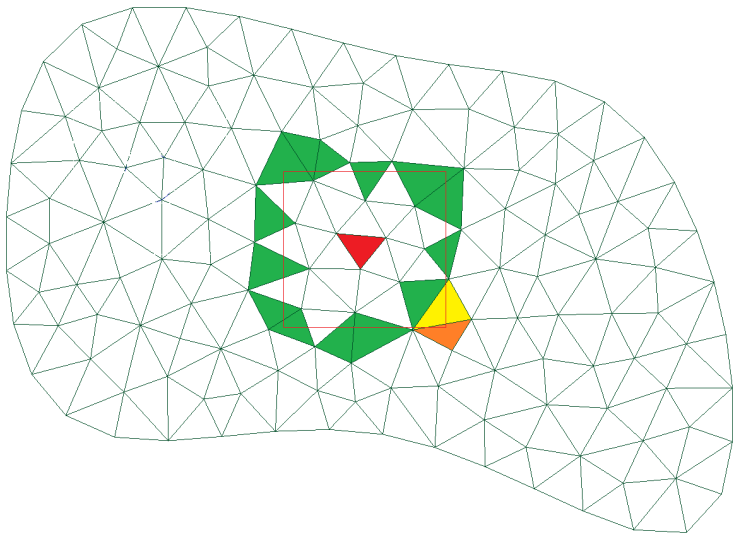
## QBX Considerations

- ▶ Global vs local approaches involve all source points, or only the "near-field" respectively
- ▶ Layer potentials provide physical meaningful off-surface potential
- ▶ Volume potentials involve arbitrary choice of "off-volume" potential

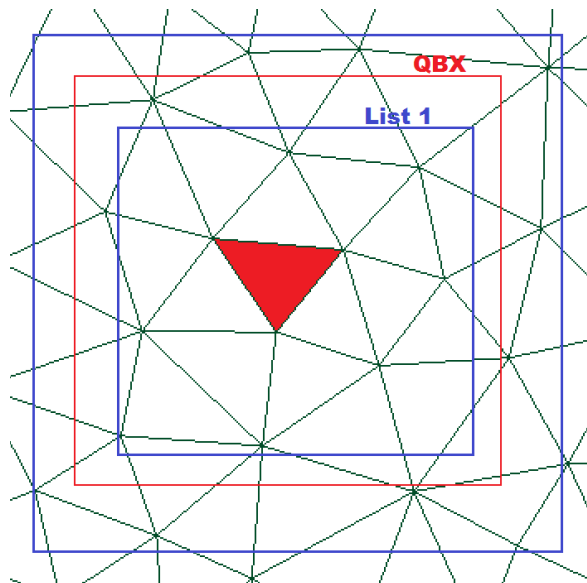
# Mesh Interaction

- ▶ Given some mesh with a spatially varying blob of "charge", local QBX needs only some of the mesh
- ▶ How are varying intersection cases to be handled?
- ▶ How conservative/efficient can one be with intersections?

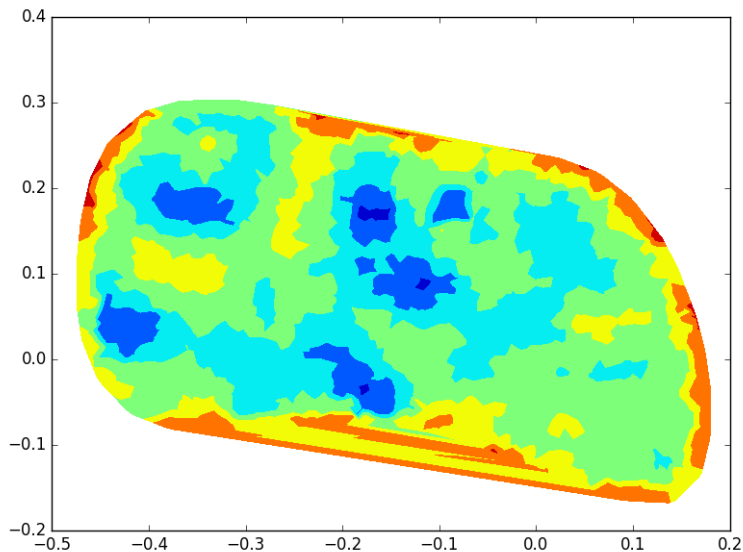
## Mesh Interaction



## FMM-QBX Interaction

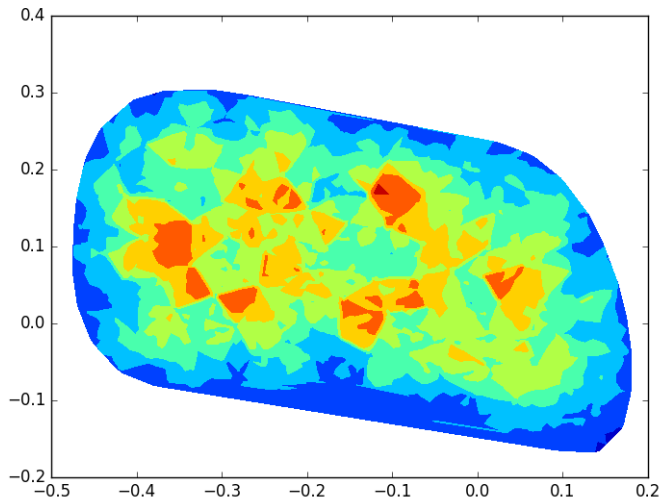


## Corrected FMM Contributions

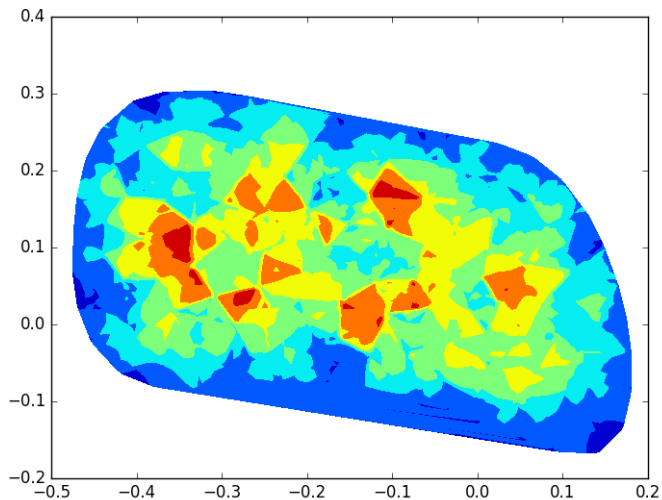




## QBX Contributions



## QBX-FMM Overlap Corrections



# Result of Combinations

