APPM4720/5720 — Homework 3

Everyone should do Problem 3.1. Then you have a choice to do either 3.2 and 3.3 or 3.4 or 3.5.

Problem 3.1: The objective of this problem is to computationally investigate the error incurred by truncating multipole expansions. Consider the following geometry: Let Ω_{τ} and Ω_{σ} be two well-separated boxes with centers c_{τ} and c_{σ} . Let c_{σ} be a source point and let c_{σ} be a target point. Consider the error function

$$e(P) = \sup \left\{ \left| \log |\boldsymbol{x} - \boldsymbol{y}| - \mathsf{B}_P(\boldsymbol{x}, \, \boldsymbol{c}_\tau) \, \mathsf{Z}_P(\boldsymbol{c}_\tau, \, \boldsymbol{c}_\sigma) \, \mathsf{C}_P(\boldsymbol{c}_\sigma, \, \boldsymbol{y}) \right| : \, \boldsymbol{x} \in \Omega_\sigma \, \, \boldsymbol{y} \in \Omega_\tau \right\}$$

where P is the length of the multipole expansion, and where

- $\mathsf{C}_P(oldsymbol{c}_\sigma,\,oldsymbol{y})\in\mathbb{C}^{(P+1)\times 1}$ maps a source to an outgoing expansion
- $\mathsf{Z}_P(\boldsymbol{c}_\tau,\,\boldsymbol{c}_\sigma)\in\mathbb{C}^{(P+1)\times(P+1)}$ maps an outgoing expansion to an incoming expansion
- $\mathsf{B}_P(\boldsymbol{x},\,\boldsymbol{c}_{\tau})\in\mathbb{C}^{1\times(P+1)}$ maps an incoming expansion to a target
- (a) Estimate e(P) experimentally for the geometry:

$$\Omega_{\sigma} = [-1, 1] \times [-1, 1], \qquad \Omega_{\tau} = [3, 5] \times [-1, 1].$$

- (b) Fit the function you determined in (a) to a curve $e(P) \sim c \cdot \alpha^P$. What is α ?
- (c) Is the supremum for a given P attained for any specific pair $\{x, y\}$? If so, find (experimentally) the pair. Does the choice depend on P?
- (d) Repeat questions (a), (b), (c) for a different geometry of your choice. (Provide a picture.)
- (e) (Optional) Can you support your observations with analysis?

Hint: The provided file main_T_ops_are_fun.m might be useful.

Problem 3.2: The objective of this exercise is to familiarize yourself with the provided prototype FMM. The questions below refer to the basic FMM provided in the file main_fmm.m when executed on a uniform particle distribution. For this case, precompute only the translation operators $\mathsf{T}^{(\text{ofo})}$, $\mathsf{T}^{(\text{ifo})}$, and $\mathsf{T}^{(\text{ifi})}$ (i.e. set flag_precomp=0).

(a) Estimate and plot the execution time of the FMM for the choices

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N_{\text{tot}} = 1\,000, \, 2\,000, \, 4\,000, \, 8\,000, \, 16\,000, \, 32\,000, \, 64\,000.
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Set nmax=50. Provide plots that track the following costs:

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t_{
m tot} total execution time, including initialization. t_{
m init} cost of initialization (computing the tree, the object T_OPS, etc.). t_{
m ofs} cost of applying T<sup>(ofs)</sup>. t_{
m ofo} cost of applying T<sup>(ofo)</sup>. t_{
m ifo} cost of applying T<sup>(ifo)</sup>. t_{
m ifo} cost of applying T<sup>(ifo)</sup>.
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 $t_{\rm tfi}$ cost of applying $\mathsf{T}^{({\rm tfi})}$.

 $t_{\rm close}$ cost of directly evaluating close range interactions.

(b) Repeat exercise (a) but now for a few different choices of nmax. Which one is the best one? Provide a new plot of the times required for this optimal choice.

Problem 3.3: Repeat Problem 3.2 but now use a non-uniform point distribution of your choice.

Problem 3.4: Can you think of a better way of computing the interaction lists? Here "better" could mean either a cleaner code that executes in more or less the same time, or a code that executes significantly faster than the provided one. If your code is **both** cleaner and faster then so much the better!

Problem 3.5: Code up the single-level Barnes-Hut method and investigate computationally how many boxes you should use for optimal performance for any given precision and given total number N_{tot} of charges. Create a plot of the best possible time t_{optimal} for several N_{tot} and estimate the dependence of t_{optimal} on N_{tot} . To keep things simple, consider only uniform particle distributions. You need only consider a fixed precision (say P = 10) but an ambitious solution should compute the optimal time for several different choices (say P = 5, 10, 15, 20).