CB\&B752/MCDB452/MB\&B752/MCDB752/CPSC752
Homework 1
Due: April 15, 2011

## Problem 1:

For the BLN model protein discussed in class and in J. D. Honeycutt and D. Thirumalai, "The nature of folded states of globular proteins," Biopolymers 32 (1992) 695, which includes Lennard-Jones-like long-range $\left(\mathrm{V}_{\mathrm{lr}}\right)$, bond length $\left(\mathrm{V}_{\mathrm{bl}}\right)$, bond angle $\left(\mathrm{V}_{\mathrm{ba}}\right)$, and dihedral angle ( $\mathrm{V}_{\mathrm{da}}$ ) interactions, calculate analytical expressions for the $\mathrm{x}-, \mathrm{y}$-, and z components of the force on a given monomer i for a chain of length $N$. The total potential energy is given by
$V_{\text {tot }}=\sum_{i=1}^{N} \sum_{j=1}^{i-2} V_{l r}\left(r_{i, j}\right)+\sum_{i=1}^{N-1} V_{b l}\left(r_{i, i+1}\right)+\sum_{i=1}^{N-2} V_{b a}\left(\theta_{i, i+1, i+2}\right)+\sum_{i=1}^{N-3} V_{d a}\left(\phi_{i, i+1, i+2, i+3}\right)$, where
$r_{i, j}=\left|\vec{r}_{i}-\vec{r}_{j}\right|, \vec{r}_{i}$ locates the center of monomer $\mathrm{i}, \sigma$ is the diameter of the monomer, $V_{l r}\left(r_{i, j}\right)=4 \varepsilon_{h}\left[\left(\frac{\sigma}{r_{i, j}}\right)^{12}+C\left(\frac{\sigma}{r_{i, j}}\right)^{6}\right]$ with $\mathrm{C}=-1$ and $\varepsilon=\varepsilon_{\mathrm{h}}$ when $\mathrm{i}, \mathrm{j}=\mathrm{B}, \mathrm{B}, \mathrm{C}=1$ and $\varepsilon=\varepsilon_{\mathrm{L}}$
$=2 / 3 \varepsilon_{\mathrm{h}}$ when $\mathrm{i}, \mathrm{j}=\mathrm{LL}, \mathrm{LB}$, and $\mathrm{C}=0$ and $\varepsilon=\varepsilon_{\mathrm{h}}$ when $\mathrm{i}, \mathrm{j}=\mathrm{NN}, \mathrm{NL}, \mathrm{NB}$, and $\varepsilon_{\mathrm{h}}$ is the LennardJones energy scale, $V_{b l}\left(r_{i, j}\right)=\frac{k_{b}}{2}\left(r_{i, j}-\sigma\right)^{2}, \mathrm{k}_{\mathrm{b}}$ is the spring constant,
$V_{b a}\left(\theta_{i, j, k}\right)=\frac{k_{\theta}}{2}\left(\theta_{i, j, k}-\theta_{0}\right)^{2}, \mathrm{k}_{\theta}$ is the bend spring constant, $\theta_{i, j, k}=\cos ^{-1}\left(\frac{\vec{r}_{i, j} \cdot \vec{r}_{k, j}}{r_{i, j} r_{k, j}}\right)$, and $\theta_{0}$ is the equilibrium bend angle, $V_{d a}\left(\phi_{i, j, k, l}\right)=A\left(1+\cos \left(\phi_{i, j, k, l}\right)\right)+B\left(1+\cos \left(3 \phi_{i, j, k, l}\right)\right)$, where A and B constants and where $\phi_{\mathrm{I}, \mathrm{j}, \mathrm{k}, 1}$ is defined by
$\cos \left(\phi_{i, j, k, l}\right)=\left(\vec{r}_{i, j} \times \vec{r}_{k, j}\right) \cdot\left(\vec{r}_{j, k} \times \vec{r}_{l, k}\right) /\left(\left|\left(\vec{r}_{i, j} \times \vec{r}_{k, j}\right) \|\left(\vec{r}_{j, k} \times \vec{r}_{l, k}\right)\right|\right)$ and
$\sin \left(\phi_{i, j, k, l}\right)=\vec{r}_{j, k} \cdot\left[\left(\vec{r}_{i, j} \times \vec{r}_{k, j}\right) \times\left(\vec{r}_{j, k} \times \vec{r}_{l, k}\right)\right] /\left(\mid\left(\vec{r}_{i, j} \times \vec{r}_{k, j}\right) \|\left(\vec{r}_{j, k} \times \vec{r}_{l, k}\right)\right)$. Use $\vec{F}_{i}=-\left[\hat{x} \frac{\partial V_{\text {tot }}}{\partial x_{i}}+\hat{y} \frac{\partial V_{\text {tot }}}{\partial y_{i}}+\hat{z} \frac{\partial V_{\text {tot }}}{\partial z_{i}}\right]$ to calculate the total force on monomer i.

## Problem 2:

Derive the following improved version of Equation (10) in J. D. Honeycutt and D. Thirumalai, "The nature of folded states of globular proteins," Biopolymers 32 (1992) 695:
$\dot{\vec{r}}_{i}(t+\Delta t)=\dot{\vec{r}}_{i}(t)\left\lfloor 1-\gamma \Delta t+\frac{(\gamma \Delta t)^{2}}{2}\right\rfloor+\Delta t\left[\left(\frac{1}{2}-\frac{\gamma \Delta t}{6}\right) \frac{\vec{F}_{i}(t)}{m_{i}}+\left(\frac{1}{2}-\frac{\gamma \Delta t}{3}\right) \frac{\vec{F}_{i}(t+\Delta t)}{m_{i}}\right]+\frac{1}{m_{i}} \int_{t}^{t+\Delta t} e^{-\gamma\left(t+\Delta t-t^{\prime}\right)} \Gamma_{i}\left(t^{\prime}\right) d t^{\prime}$

## Problem 3 (CPSC752):

Write a c++ function to calculate the N-3 dihedral angles for a polymer chain, assuming that the positions $\vec{r}_{i}$ of the $\mathrm{i}=1, \mathrm{~N}$ monomers are specified. Download the file positions.dat from the course website, which lists the positions of the 46 monomers corresponding to the ground state of the BLN model $\mathrm{B}_{9} \mathrm{~N}_{3}(\mathrm{LB})_{4} \mathrm{~N}_{3} \mathrm{~B}_{9} \mathrm{~N}_{3}(\mathrm{LB})_{5} \mathrm{~L}$ in the following format:
$\mathrm{r}_{\mathrm{x} 1} \mathrm{r}_{\mathrm{y} 1} \mathrm{r}_{\mathrm{z} 1}$
$\mathrm{r}_{\mathrm{x} 2} \mathrm{r}_{\mathrm{y} 2} \mathrm{r}_{\mathrm{z} 2}$
...
$\mathrm{r}_{\mathrm{xN}} \mathrm{r}_{\mathrm{yN}} \mathrm{r}_{\mathrm{zN}}$
Compare your results to the answers in file dihedral.dat, which can also be downloaded from course website.

