

- We are doing md simulation with PBC of a 2d system. We try to verify our result using various benchmark. Configuration data is given in file: config_benchmark3.txt

Antik da Result:

```
density = 1.2
N = 100
r_c = 2.5
particle type A
```

```
double force_check = 0.e0;
for(int i=0;i<N;i++) force_check+=(sphere[i].Fx*sphere[i].Fx)+(sphere[i].Fy*sphere[i].Fy);
force_check = sqrt(force_check);
printf("%.16lf\n",force_check);
```

```
force_check = 0.0000000036204816
```

My Result:

```
Arrays are allocated.
-----
Initial force given to system.
-----
Force Disbalance: 0.0000000036204378
arnab@arnab-phys:~/Project_dpl/NVE_PBC_2d_basic$
```

```
arnab@arnab-phys:~/Project_dpl/NVE_PBC_2d_basic$ gcc md_p
arnab@arnab-phys:~/Project_dpl/NVE_PBC_2d_basic$ ./a.out
Arrays are allocated.
-----
Initial force given to system.
-----
Force Disbalance: 0.0000000036203637
arnab@arnab-phys:~/Project_dpl/NVE_PBC_2d_basic$
```

- dt check has done in test_variance.py file.

