

```
...float coory = gridspacing * yindex;
float coorx = gridspacing * xindex;
float gridspacing_coalesce = gridspacing * BLOCKSIZEX;
int atomid;
for (atomid=0; atomid<numatoms; atomid++) {
    float dy = coory - atominfo[atomid].y;
    float dyz2 = (dy * dy) + atominfo[atomid].z;
    float dx1 = coorx - atominfo[atomid].x;
[...]
```

float dx8 = dx7 + gridspacing_coalesce;

```
energyvalx1 += atominfo[atomid].w * rsqrtf(dx1*dx1 + dyz2);
[...]
```

energyvalx8 += atominfo[atomid].w * rsqrtf(dx8*dx8 + dyz2);

```
}
energygrid[outaddr
[...]
```

energygrid[outaddr+7*BLOCKSIZEX] += energyvalx7;

Points spaced for
memory coalescing

Reuse partial distance
components $dy^2 + dz^2$

Global memory ops
occur only at the end
of the kernel,
decreases register use