

Surface Interaction

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Algorithm 1 Surface Interaction

```
1: procedure INTERACTIONS(lattice site,atom,n,E)
2:   for each atom make a list of its neighbors
3:   [(n=0|central atom each time, type(0)),
4:    (n=1|1st neighbor, type(1)),
5:    .....
6:    (n=5|5th neighbor, type(5))] where type is 1 or 2 depends by the number of species we have.
7:   if type(0) = type(1:5) then
8:      $E = V1$ : interaction  $V_{central-atom,j}$ 
9:   else
10:     $E = V2$ : interaction  $V_{central-atom,j}$ 
```

