Molecular Origami of BrICl^-

given information

| | | |
|---|---|---|
| ElementNames | [ (I) (Br) (Cl) ] | |
| dotted | F | |
| distance | 264.792 | $I^1$-$Br^1$ |
| distance | 265.066 | $I^1$-$Cl^1$ |
| angle | 179.605 | $Cl^1$-$I^1$-$Br^1$ |
| | 529.9 | $Cl^1$-$Br^1$ |
| dopage | T | |
| AutoAlign | F | |
| showboth | F | |

structure type: XAB

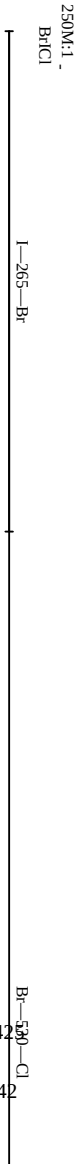
Molecular Origami of BrICl^-

!I1
Br1
Cl1
BrICl^-

linear

scale 250,000,000 : 1
units: pm
offsetx -0.07  offsety 0.9

View -1




Current:  (centerx 4.23) (centery 5.90) (scale 250)
%%BoundingBox: 295 -148 315 435  actual: 305 -138 305 425  center: 305 143  actual size: 0 563
Better:  (centerx 4.25) (centery 9.41) (scale 250)
%%BoundingBox: 291 169 311 752  actual: 301 179 301 742  center: 301 461  actual size: 0 563