Molecular Origami of CN^-

given information

| | | |
|---|---|---|
| ElementNames | [ (C) (N) ] | |
| distance | 124.491 | $C^{1}$-$N^{1}$ |
| dopage | T | |
| AutoAlign | F | |
| showboth | F | |

structure type: XA

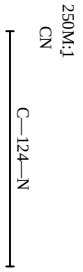
Molecular Origami of CN^-

!C1
N1
CN^-

diatomic

scale 250,000,000 : 1
units: pm
offsetx -0.05  offsety 0.9

View -1




Current:  (centerx 4.25) (centery 5.90) (scale 250)
%%BoundingBox: 296 327 316 435  actual: 306 337 306 425  center: 306 381  actual size: 0 88
Better:  (centerx 4.25) (centery 6.11) (scale 250)
%%BoundingBox: 292 407 312 515  actual: 302 417 302 505  center: 302 461  actual size: 0 88