Molecular Origami of NO

given information

| | | |
|---|---|---|
| ElementNames | [ (N) (O) ] | |
| distance | 117.567 | $N^{1}$-$O^{1}$ |
| dopage | T | |
| AutoAlign | F | |
| showboth | F | |

structure type: XA

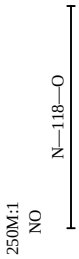
Molecular Origami of NO

!N1
O1
NO

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: 296 415 316 518  actual: 306 425 306 508  center: 306 466  actual size: 0 83
Better:  (centerx 4.25) (centery 4.92) (scale 250)
%%BoundingBox: 292 409 312 512  actual: 302 419 302 502  center: 302 461  actual size: 0 83