

# Molecular Origami of N2

given information

ElementNames	[ (N) (N) ]	
distance	110.000	N <sup>1</sup> -N <sup>2</sup>
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XA

# Molecular Origami of N2

!N1

N2

N2

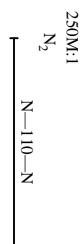
diatomic

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: 296 337 316 435 actual: 306 347 306 425

center: 306 386

actual size: 0 78

Better: (centerx 4.25) (centery 6.04) (scale 250)

%%BoundingBox: 292 412 312 510 actual: 302 422 302 500

center: 302 461

actual size: 0 78