

Molecular Origami of CN⁻

given information

ElementNames	[(C) (N)]	
distance	125.240	C ¹ -N ¹
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XA

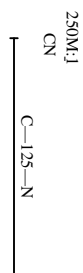
Molecular Origami of CN⁻

!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: 296 326 316 435 actual: 306 336 306 425

center: 306 380

actual size: 0 89

Better: (centerx 4.25) (centery 6.12) (scale 250)

%%BoundingBox: 292 406 312 515

actual: 302 416 302 505

center: 302 461

actual size: 0 89