Molecular Origami of N3^-

given information

ElementNames	[(N) (N) (N)]	
dotted	F	
distance	117.990	N^2-N^3
distance	118.002	N^2-N^1
angle	179.997	N^3-N^2-N^1
	236.	N^3-N^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

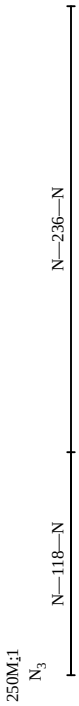
Molecular Origami of N3^-

!N2
N1
N3
N3^-

linear

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: 296 415 316 686 actual: 306 425 306 676 center: 306 550 actual size: 0 251
Better: (centerx 4.25) (centery 3.76) (scale 250)
%%BoundingBox: 292 325 312 596 actual: 302 335 302 586 center: 302 461 actual size: 0 251