Molecular Origami of NCO^-

given information

ElementNames	[(C) (N) (O)]	
dotted	F	
distance	112.730	C^1-O^1
distance	121.992	C^1-N^1
angle	179.062	O^1-C^1-N^1
	234.7	O^1-N^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

Molecular Origami of NCO^-

!C1
N1
O1
NCO^-

linear

scale 250,000,000 : 1
units: pm
offsetx -0.07 offsety 0.9

View -1

Current: (centerx 4.23) (centery 5.90) (scale 250)
%%BoundingBox: 295 162 315 435 actual: 305 172 305 425 center: 305 298 actual size: 0 253
Better: (centerx 4.25) (centery 7.26) (scale 250)
%%BoundingBox: 291 324 311 597 actual: 301 334 301 587 center: 301 461 actual size: 0 253