Molecular Origami of NO2

given information

ElementNames	[(N) (O) (O)]	
dotted	F	
distance	117.570	N^1-O^2
distance	132.501	N^1-O^1
angle	120.001	O^2-N^1-O^1
	216.7	O^2-O^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

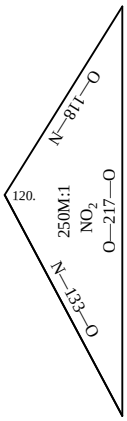
Molecular Origami of NO2

!N1
O1
O2
NO2

bent

scale 250,000,000 : 1
units: pm
offsetx 0.91 offsety 0.9

View -1



Current: (centerx 5.21) (centery 5.90) (scale 250)
%%BoundingBox: 321 338 385 512 actual: 331 348 375 502 center: 353 425 actual size: 44 154
Better: (centerx 4.56) (centery 5.50) (scale 250)
%%BoundingBox: 339 374 404 548 actual: 349 384 394 538 center: 372 461 actual size: 44 154