Molecular Origami of NO2^-

given information

ElementNames	[(N) (O) (O)]	
dotted	F	
distance	116.176	N^1-O^1
distance	126.200	N^1-O^2
angle	107.680	O^2-N^1-O^1
	195.8	O^2-O^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

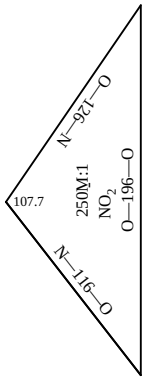
Molecular Origami of NO2^-

!N1
O1
O2
NO2^-

bent

scale 250,000,000 : 1
units: pm
offsetx 0.85 offsety 0.9

View -1



Current: (centerx 5.15) (centery 5.90) (scale 250)
%%BoundingBox: 310 345 381 504 actual: 320 355 371 494 center: 346 425 actual size: 51 139
Better: (centerx 4.60) (centery 5.50) (scale 250)
%%BoundingBox: 332 381 402 540 actual: 342 391 392 530 center: 367 461 actual size: 51 139