Molecular Origami of NO2^-

given information

ElementNames	[(N) (O) (O)]	
dotted	F	
distance	122.313	N^1-O^1
distance	122.902	N^1-O^2
angle	116.298	O^2-N^1-O^1
	208.3	O^2-O^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

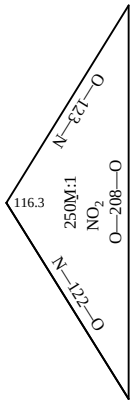
Molecular Origami of NO2^-

!N1
O1
O2
NO2^-

bent

scale 250,000,000 : 1
units: pm
offsetx 0.87 offsety 0.9

View -1



Current: (centerx 5.17) (centery 5.90) (scale 250)
%%BoundingBox: 316 341 382 509 actual: 326 351 372 499 center: 349 425 actual size: 46 148
Better: (centerx 4.57) (centery 5.50) (scale 250)
%%BoundingBox: 336 377 402 545 actual: 346 387 392 535 center: 369 461 actual size: 46 148