Molecular Origami of SO2

given information

ElementNames	[(S) (O) (O)]	
dotted	F	
distance	142.994	S^1-O^2
distance	143.002	S^1-O^1
angle	119.504	O^2-S^1-O^1
	247.1	O^2-O^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

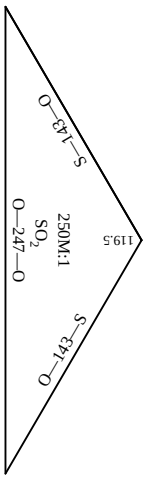
Molecular Origami of SO2

!S1
O1
O2
SO2

bent

scale 250,000,000 : 1
units: pm
offsetx -1.11 offsety 0.9

View -1



Current: (centerx 3.19) (centery 5.90) (scale 250)
%%BoundingBox: 220 327 291 522 actual: 230 337 281 512 center: 255 425 actual size: 51 175
Better: (centerx 3.90) (centery 5.50) (scale 250)
%%BoundingBox: 191 363 262 558 actual: 201 373 252 548 center: 226 461 actual size: 51 175