Molecular Origami of ClO2^-

given information

ElementNames	[(Cl) (O) (O)]	
dotted	F	
distance	165.700	Cl^1-O^1
distance	165.702	Cl^1-O^2
angle	109.498	O^2-Cl^1-O^1
	270.6	O^2-O^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

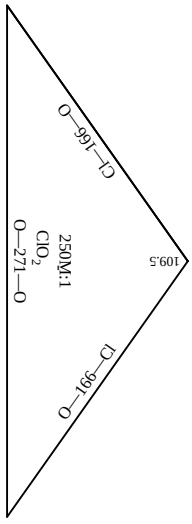
Molecular Origami of ClO2^-

!Cl1
O1
O2
ClO2^-

bent

scale 250,000,000 : 1
units: pm
offsetx -1.31 offsety 0.9

View -1



Current: (centerx 2.99) (centery 5.90) (scale 250)
%%BoundingBox: 205 319 293 531 actual: 215 329 283 521 center: 249 425 actual size: 68 192
Better: (centerx 3.78) (centery 5.50) (scale 250)
%%BoundingBox: 168 355 256 567 actual: 178 365 246 557 center: 212 461 actual size: 68 192