Molecular Origami of H2O2

given information

ElementNames	[(O) (H) (O)]	
dotted	F	
distance	99.009	O^1-H^1
distance	146.998	O^1-O^2
angle	98.003	O^2-O^1-H^1
	188.3	O^2-H^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

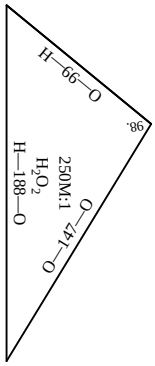
Molecular Origami of H2O2

!O1
H1
O2
H2O2

bent

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

View -1



Current: (centerx 3.43) (centery 5.90) (scale 250)
%%BoundingBox: 237 348 311 502 actual: 247 358 301 492 center: 274 425 actual size: 54 133
Better: (centerx 3.87) (centery 5.50) (scale 250)
%%BoundingBox: 206 384 280 538 actual: 216 394 270 528 center: 243 461 actual size: 54 133