Molecular Origami of H2O

given information

ElementNames	[(O) (H) (H)]	
dotted	F	
distance	93.668	O^1-H^1
distance	95.273	O^1-H^2
angle	109.110	H^2-O^1-H^1
	153.9	H^2-H^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

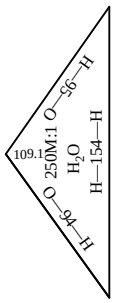
Molecular Origami of H2O

!O1
H1
H2
H2O

bent

scale 250,000,000 : 1
units: pm
offsetx 0.66 offsety 0.9

View -1



Current: (centerx 4.96) (centery 5.90) (scale 250)
%%BoundingBox: 308 360 367 489 actual: 318 370 357 479 center: 338 425 actual size: 39 109
Better: (centerx 4.52) (centery 5.50) (scale 250)
%%BoundingBox: 324 396 383 525 actual: 334 406 373 515 center: 354 461 actual size: 39 109