Molecular Origami of H2O

given information

ElementNames	[(O) (H) (H)]	
dotted	F	
distance	89.413	O^1-H^1
distance	93.942	O^1-H^2
angle	111.667	H^2-O^1-H^1
	151.7	H^2-H^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

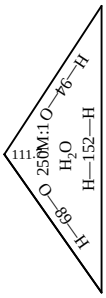
Molecular Origami of H2O

!O1
H1
H2
H2O

bent

scale 250,000,000 : 1
units: pm
offsetx 0.63 offsety 0.9

View -1



Current: (centerx 4.93) (centery 5.90) (scale 250)
%%BoundingBox: 309 361 365 489 actual: 319 371 355 479 center: 337 425 actual size: 36 108
Better: (centerx 4.50) (centery 5.50) (scale 250)
%%BoundingBox: 323 397 380 525 actual: 333 407 370 515 center: 351 461 actual size: 36 108