

Molecular Origami of H2O

given information

ElementNames	[(O) (H) (H)]
dotted	F
distance	93.078
distance	97.142
angle	108.160
	154.1
dopage	T
AutoAlign	F
showboth	F

O¹-H²
O¹-H¹
H²-O¹-H¹
H²-H¹

structure type: XAB

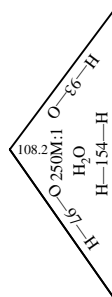
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!O1
H1
H2
H2O

bent

scale 250,000,000 : 1
units: pm
offsetx 0.68 offsety 0.9

View -1



Current: (centerx 4.98) (centery 5.90) (scale 250)

%%BoundingBox: 309 360 369 489 actual: 319 370 359 479

Better: (centerx 4.52) (centery 5.50) (scale 250)

%%BoundingBox: 325 396 385 525 actual: 335 406 375 515

center: 339 425

actual size: 40 109

center: 355 461

actual size: 40 109