Molecular Origami of HOBr

given information

ElementNames	[(O) (Br) (H)]	
dotted	F	
distance	95.327	O^1-H^1
distance	184.999	O^1-Br^1
angle	109.998	H^1-O^1-Br^1
	235.3	H^1-Br^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

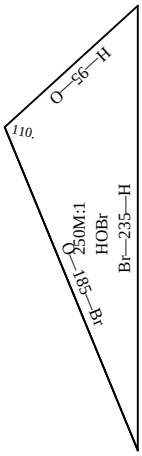
Molecular Origami of HOBr

!O1
Br1
H1
HOBr

bent

scale 250,000,000 : 1
units: pm
offsetx 1.1 offsety 1.01

View -1



Current: (centerx 5.40) (centery 6.01) (scale 250)
%%BoundingBox: 329 339 399 526 actual: 339 349 389 516 center: 364 433 actual size: 50 167
Better: (centerx 4.60) (centery 5.50) (scale 250)
%%BoundingBox: 350 375 420 562 actual: 360 385 410 552 center: 385 469 actual size: 50 167