Molecular Origami of HOF

given information

ElementNames	[(O) (F) (H)]	
dotted	F	
distance	95.316	O^1-H^1
distance	144.007	O^1-F^1
angle	96.996	H^1-O^1-F^1
	182.1	H^1-F^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

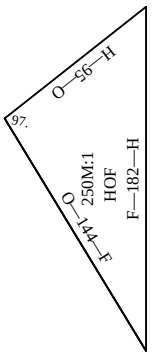
Molecular Origami of HOF

!O1
F1
H1
HOF

bent

scale 250,000,000 : 1
units: pm
offsetx 0.82 offsety 1.11

View -1



Current: (centerx 5.12) (centery 6.11) (scale 250)
%%BoundingBox: 306 365 379 514 actual: 316 375 369 504 center: 342 440 actual size: 53 129
Better: (centerx 4.62) (centery 5.50) (scale 250)
%%BoundingBox: 329 401 402 550 actual: 339 411 392 540 center: 365 476 actual size: 53 129