Molecular Origami of HOCl

given information

ElementNames	[(O) (Cl) (H)]	
dotted	F	
distance	95.335	O^1-H^1
distance	169.985	O^1-Cl^1
angle	113.003	H^1-O^1-Cl^1
	225.1	H^1-Cl^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

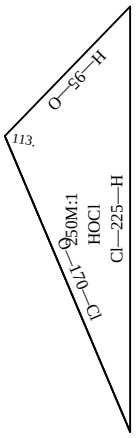
Molecular Origami of HOCl

!O1
Cl1
H1
HOCl

bent

scale 250,000,000 : 1
units: pm
offsetx 1.06 offsety 0.9

View -1



Current: (centerx 5.36) (centery 5.90) (scale 250)
%%BoundingBox: 329 335 396 515 actual: 339 345 386 505 center: 362 425 actual size: 47 159
Better: (centerx 4.58) (centery 5.50) (scale 250)
%%BoundingBox: 349 371 416 551 actual: 359 381 406 541 center: 382 461 actual size: 47 159