Molecular Origami of Cu(PPh3)3^+

given information

ElementNames	[(Cu) (P) (P) (P)]	
dotted	F	
distance	229.289	Cu^1-P^1
distance	229.565	Cu^1-P^3
distance	229.667	Cu^1-P^2
angle	117.464	P^2-Cu^1-P^1
	392.3	P^2-P^1
angle	119.965	P^3-Cu^1-P^1
	397.3	P^3-P^1
angle	122.569	P^3-Cu^1-P^2
	402.8	P^3-P^2
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

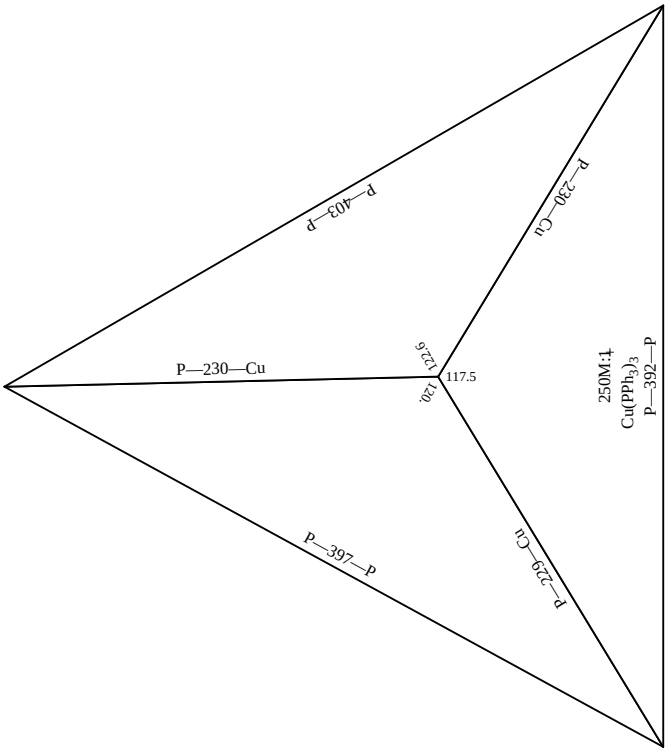
Molecular Origami of Cu(PPh3)3^+

!Cu1
P1
P2
P3
Cu(PPh3)3^+

trigonal planar

scale 250,000,000 : 1
units: pm
offsetx 1.67 offsety 0.9

View -1



Current: (centerx 5.97) (centery 5.90) (scale 250)
%%BoundingBox: 173 276 440 574 actual: 183 286 430 564 center: 306 425 actual size: 247 278
Better: (centerx 5.97) (centery 5.50) (scale 250)
%%BoundingBox: 293 312 560 610 actual: 303 322 550 600 center: 426 461 actual size: 247 278