Molecular Origami of CH2I2

given information

ElementNames	[(C) (H) (H) (I) (I)]	
distance	108.663	C^1-H^2
distance	110.399	C^1-H^1
distance	214.323	C^1-I^2
distance	215.801	C^1-I^1
angle	106.698	I^2-C^1-H^2
	266.7	I^2-H^2
angle	107.421	I^1-C^1-H^1
	270.2	I^1-H^1
angle	109.157	I^1-C^1-H^2
	271.6	I^1-H^2
angle	109.928	H^2-C^1-H^1
	179.4	H^2-H^1
angle	110.759	I^2-C^1-H^1
	273.7	I^2-H^1
angle	112.868	I^2-C^1-I^1
	358.4	I^2-I^1
dopage	T	
AutoAlign	F	

structure type: XABCD

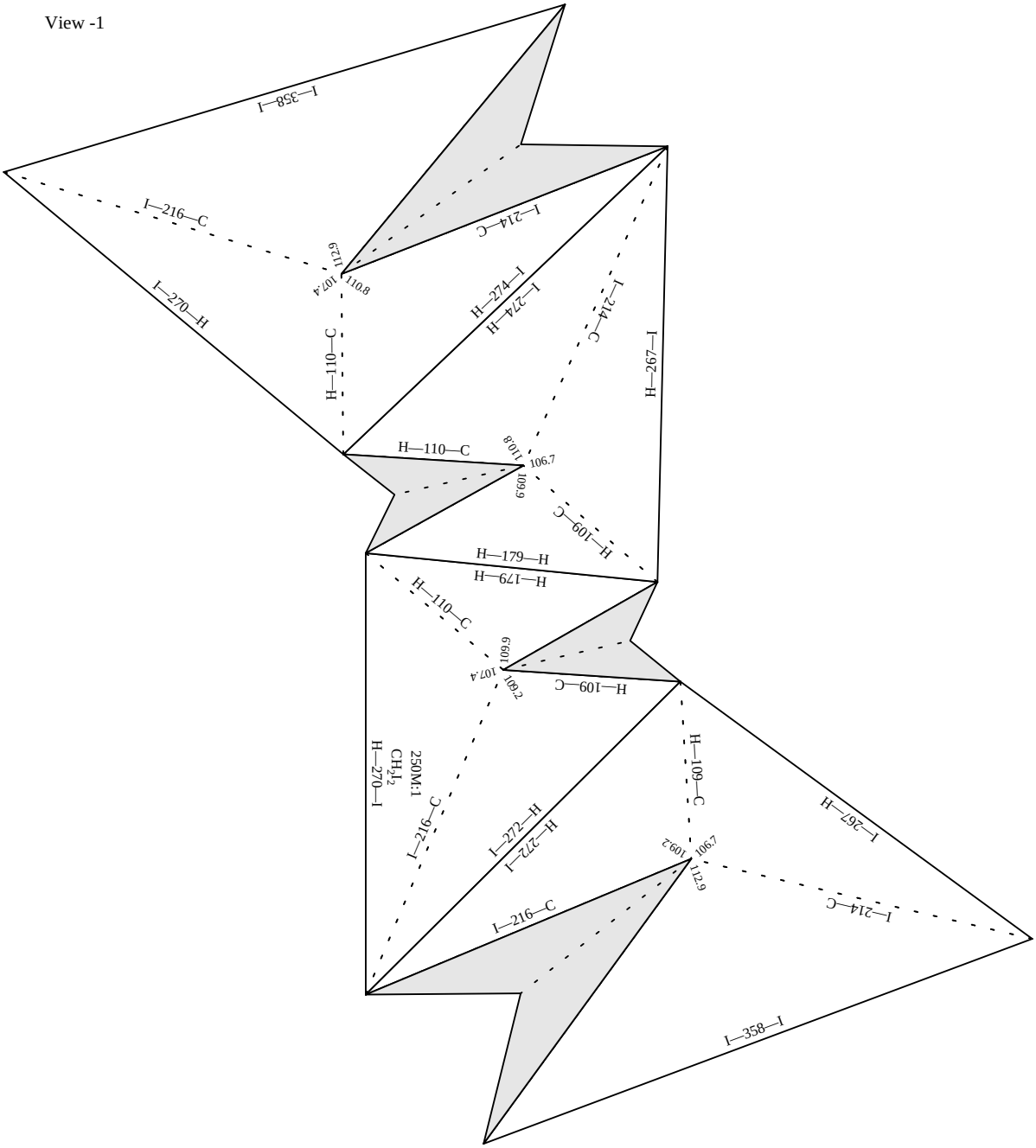
Molecular Origami of CH2I2

!C1
H1
H2
I1
I2
CH2I2

special tetrahedral

scale 250,000,000 : 1
units: pm
offsetx -0.54 offsety -0.63

View -1



Current: (centerx 3.76) (centery 4.37) (scale 250)
%%BoundingBox: 44 99 510 613 actual: 54 109 500 603 center: 277 356 actual size: 446 494
Better: (centerx 4.16) (centery 4.92) (scale 250)
%%BoundingBox: 34 94 500 608 actual: 44 104 490 598 center: 267 351 actual size: 446 494