Molecular Origami of CH3I

given information

ElementNames	[(C) (H) (H) (H) (I)]	
distance	108.251	C^1-H^3
distance	108.259	C^1-H^1
distance	111.042	C^1-H^2
distance	207.738	C^1-I^1
angle	105.871	H^2-C^1-H^1
	175.	H^2-H^1
angle	105.873	H^3-C^1-H^2
	175.	H^3-H^2
angle	110.209	I^1-C^1-H^2
	267.3	I^1-H^2
angle	111.467	H^3-C^1-H^1
	178.9	H^3-H^1
angle	111.554	I^1-C^1-H^1
	267.2	I^1-H^1
angle	111.556	I^1-C^1-H^3
	267.2	I^1-H^3
dopage	T	
AutoAlign	F	

structure type: XABCD

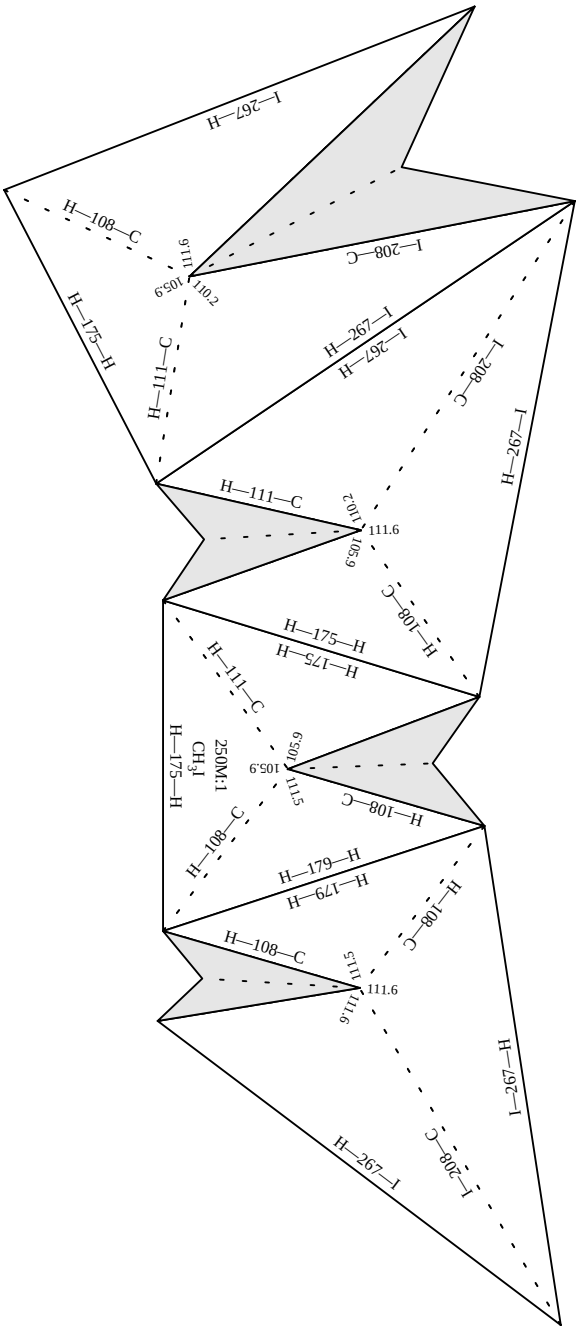
Molecular Origami of CH3I

!C1
H1
H2
H3
I1
CH3I

special tetrahedral

scale 250,000,000 : 1
units: pm
offsetx -0.04 offsety 0.38

View -1



Current: (centerx 4.26) (centery 5.38) (scale 250)
%%BoundingBox: 190 169 424 683 actual: 200 179 414 673 center: 307 426 actual size: 214 494
Better: (centerx 4.24) (centery 4.96) (scale 250)
%%BoundingBox: 186 166 420 681 actual: 196 176 410 671 center: 303 423 actual size: 214 494