Molecular Origami of CH3I

given information

ElementNames	[(C) (H) (H) (H) (I)]	
distance	99.992	C^1-H^1
distance	99.999	C^1-H^3
distance	100.000	C^1-H^2
distance	209.971	C^1-I^1
angle	109.439	H^3-C^1-H^1
	163.3	H^3-H^1
angle	109.439	H^2-C^1-H^1
	163.3	H^2-H^1
angle	109.448	H^3-C^1-H^2
	163.3	H^3-H^2
angle	109.500	I^1-C^1-H^3
	261.	I^1-H^3
angle	109.500	I^1-C^1-H^1
	261.	I^1-H^1
angle	109.502	I^1-C^1-H^2
	261.	I^1-H^2
dopage	T	
AutoAlign	F	

structure type: XABCD

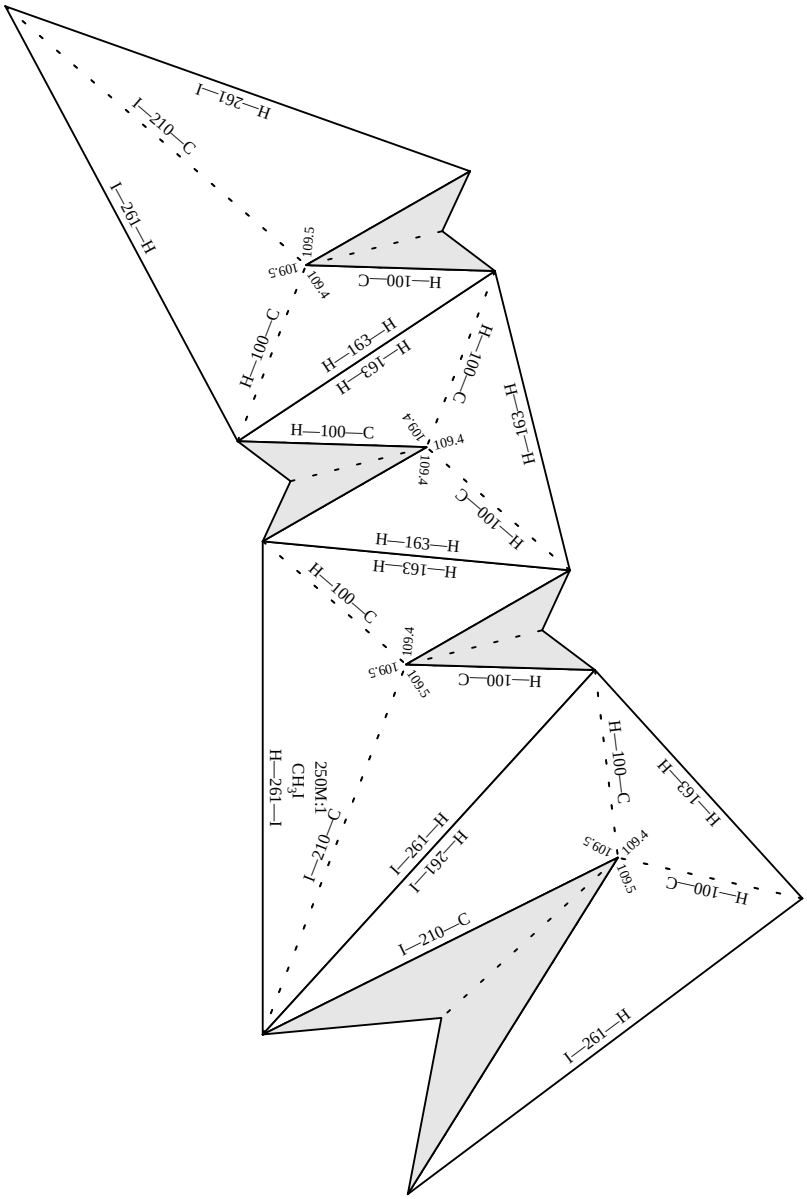
Molecular Origami of CH3I

!C1
H1
H2
H3
I1
CH3I

special tetrahedral

scale 250,000,000 : 1
units: pm
offsetx -0.43 offsety -0.44

View -1



Current: (centerx 3.87) (centery 4.56) (scale 250)
%%BoundingBox: 118 120 437 585 actual: 128 130 427 575 center: 278 352 actual size: 299 445
Better: (centerx 4.26) (centery 5.17) (scale 250)
%%BoundingBox: 116 132 434 597 actual: 126 142 424 587 center: 275 364 actual size: 299 445