

Molecular Origami of H2S

given information

ElementNames	[(S) (H) (H)]
dotted	F
distance	135.000
distance	135.085
angle	92.036
	194.3
dopage	T
AutoAlign	F
showboth	F

S¹-H¹
S¹-H²
H²-S¹-H¹
H²-H¹

structure type: XAB

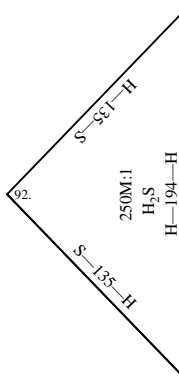
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!S1
H1
H2
H2S

bent

scale 250,000,000 : 1
units: pm
offsetx 0.91 offsety 0.9

View -1



Current: (centerx 5.21) (centery 5.90) (scale 250)

%%BoundingBox: 299 346 385 504 actual: 309 356 375 494

Better: (centerx 4.71) (centery 5.50) (scale 250)

%%BoundingBox: 328 382 415 540 actual: 338 392 405 530

center: 342 425

actual size: 66 138

center: 372 461

actual size: 66 138