

Molecular Origami of SF2

given information

ElementNames	[(S) (F) (F)]	
dotted	F	
distance	159.000	S ¹ -F ¹
distance	159.043	S ¹ -F ²
angle	97.987	F ² -S ¹ -F ¹
	240.	F ² -F ¹
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

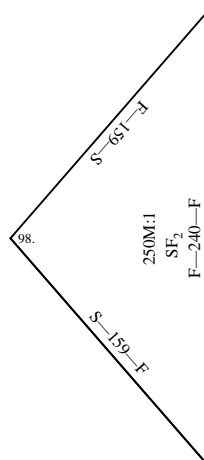
Molecular Origami of SF2

!S1
F1
F2
SF2

bent

scale 250,000,000 : 1
units: pm
offsetx 1.12 offsety 0.9

View -1



Current: (centerx 5.42) (centery 5.90) (scale 250)

%%BoundingBox: 306 330 400 520 actual: 316 340 390 510

center: 353 425

actual size: 74 170

Better: (centerx 4.76) (centery 5.50) (scale 250)

%%BoundingBox: 340 366 434 556 actual: 350 376 424 546

center: 387 461

actual size: 74 170