2-[[4-chloro-6-(cyclopropylamino)-1,3,5-triazin-2-y

Inchi: InChI=1S/C10H13ClN6/c1-10(2,5-12)17-9-15-7(11)14-8(16-9)13-6-3-4-6/h6H,3-4H2,1-2H
InchiKey: WUZNHSBFPPFULJ-UHFFFAOYSA-N
Formula: C10H13ClN6
SMILES: CC(C)(C#N)Nc1nc(Cl)nc(NC2CC2)n1
Mol. weight [g/mol]: 252.71

Physical Properties

Property code	Value	Unit	Source
logp	1.813		Crippen Method
mcvol	180.660	ml/mol	McGowan Method
tf	430.37	K	Cheméo Relay (1.0)

Sources

McGowan Method: http://link.springer.com/article/10.1007/BF02311772
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
Cheméo Relay (1.0):
Cheméo Relay (1.0, beta):
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=C32889488&Units=SI

Legend

logp: Octanol/Water partition coefficient
mcvol: McGowan's characteristic volume
tf: Normal melting (fusion) point