

6-Chloro-N,N'-(1,3-dimethylbutyl)-[1,3,5]triazine-2,

Inchi: InChI=1S/C15H28ClN5/c1-9(2)7-11(5)17-14-19-13(16)20-15(21-14)18-12(6)8-10(3)4/h9-
InchiKey: HSYFFQOHBSEBHZ-UHFFFAOYSA-N
Formula: C15H28ClN5
SMILES: CC(C)CC(C)Nc1nc(Cl)nc(NC(C)CC(C)C)n1
Mol. weight [g/mol]: 313.87

Physical Properties

Property code	Value	Unit	Source
log10ws	-5.40		Crippen Method
logp	4.218		Crippen Method
mcpvol	260.590	ml/mol	McGowan Method
rinpol	2123.87		NIST Webbook
rinpol	2188.26		NIST Webbook
rinpol	2123.87		NIST Webbook
rinpol	2138.81		NIST Webbook
rinpol	2151.76		NIST Webbook
rinpol	2188.26		NIST Webbook
rinpol	2203.23		NIST Webbook
rinpol	2123.87		NIST Webbook
rinpol	2188.26		NIST Webbook

Sources

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
McGowan Method: <http://link.springer.com/article/10.1007/BF02311772>
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R288647&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>

Legend

log10ws: Log10 of Water solubility in mol/l

logp: Octanol/Water partition coefficient
mcvol: McGowan's characteristic volume
rinpol: Non-polar retention indices

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