N<1>,n<2>-bis[4-(1-aziridinyl)butyl]ethanediamide

Other names: N<1>,n<2>-bis[4-(1-aziridinyl)butyl]ethanediamide
Inchi: InChI=1S/C14H26N4O2/c19-13(15-5-1-3-7-17-9-10-17)14(20)16-6-2-4-8-18-11-12-18/h1
InchiKey: RJZZFXJZXSMGDS-UHFFFAOYSA-N
Formula: C14H26N4O2
SMILES: O=C(NCCCCN1CC1)C(=O)NCCCCN1CC1
Mol. weight [g/mol]: 282.39

Physical Properties

Property code	Value	Unit	Source
logp	-0.590		Crippen Method
mcvol	229.460	ml/mol	McGowan Method

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=C24480260&Units=SI

Legend

logp: Octanol/Water partition coefficient
mcvol: McGowan's characteristic volume