Glutarimide, N-(4-methoxybenzyl)-

Inchi: InChI=1S/C13H15NO3/c1-17-11-7-5-10(6-8-11)9-14-12(15)3-2-4-13(14)16/h5-8H,2-4,9H
InchiKey: OIGPWDHSJVVQHY-UHFFFAOYSA-N
Formula: C13H15NO3
SMILES: COc1ccc(CN2C(=O)CCCC2=O)cc1
Mol. weight [g/mol]: 233.26

Physical Properties

Property code	Value	Unit	Source
log10ws	-2.58		Crippen Method
logp	1.734		Crippen Method
mcvol	178.400	ml/mol	McGowan Method
rinpol	2153.00		NIST Webbook

Sources

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=U360201&Units=SI
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
McGowan Method: http://link.springer.com/article/10.1007/BF02311772

Legend

log10ws: Log10 of Water solubility in mol/l
logp: Octanol/Water partition coefficient
mcvol: McGowan's characteristic volume
rinpol: Non-polar retention indices