

Californine-M, (di-(demethylene-methyl-)), isomer-3, 2AC

Inchi: InChI=1S/C23H25NO6/c1-12(25)29-22-9-15-7-18-16-10-21(28-5)20(27-4)8-14(16)6-19(2)
InchiKey: FZHLVTLLKLWVSS-UHFFFAOYSA-N
Formula: C23H25NO6
SMILES: COc1cc2c(cc1OC)C1Cc3cc(OC(C)=O)c(OC(C)=O)cc3C(C2)N1C
Mol. weight [g/mol]: 411.45

Physical Properties

Property code	Value	Unit	Source
log10ws	-5.22		Crippen Method
logp	3.381		Crippen Method
mcvol	302.290	ml/mol	McGowan Method
rinpol	3055.00		NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R289011&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

Latest version available from:

<https://www.chemeo.com/cid/25-527-9/Californine-M-di-demethylene-methyl-isomer-3-2AC.pdf>

Generated by Cheméo on 2025-12-05 21:53:58.845056479 +0000 UTC m=+4719836.375097132.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.