

2-ethyl-4-propylthiazole

Inchi: InChI=1S/C8H13NS/c1-3-5-7-6-10-8(4-2)9-7/h6H,3-5H2,1-2H3
InchiKey: DDNSNWZYKNNLRJ-UHFFFAOYSA-N
Formula: C8H13NS
SMILES: CCCc1csc(CC)n1
Mol. weight [g/mol]: 155.26

Physical Properties

Property code	Value	Unit	Source
log10ws	-3.02		Crippen Method
logp	2.658		Crippen Method
mcvol	130.450	ml/mol	McGowan Method
ripol	1528.00		NIST Webbook
ripol	1528.00		NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R298936&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
Crippen Method: https://www.cheméo.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
ripol: Polar retention indices

Latest version available from:

<https://www.cheméo.com/cid/81-381-9/2-ethyl-4-propylthiazole.pdf>

Generated by Cheméo on 2024-04-25 05:39:23.478513315 +0000 UTC m=+16312812.399090630.

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