

Thiophene-2-carboxamide, N-ethyl-N-tetradecyl-

Inchi: InChI=1S/C21H37NOS/c1-3-5-6-7-8-9-10-11-12-13-14-15-18-22(4-2)21(23)20-17-16-19-
InchiKey: XPWKAJYQYHMBHF-UHFFFAOYSA-N
Formula: C21H37NOS
SMILES: CCCCCCCCCCCCCCN(CC)C(=O)c1cccs1
Mol. weight [g/mol]: 351.59

Physical Properties

Property code	Value	Unit	Source
log10ws	-7.23		Crippen Method
logp	6.911		Crippen Method
mcvol	315.190	ml/mol	McGowan Method
rinpol	3166.00		NIST Webbook
rinpol	3166.00		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=U415316&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>

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/94-655-1/Thiophene-2-carboxamide-N-ethyl-N-tetradecyl.pdf>

Generated by Cheméo on 2024-04-17 15:25:55.454216398 +0000 UTC m=+15656804.374793714.

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