D-Proline, N-ethoxycarbonyl-, nonyl ester

Inchi: InChI=1S/C17H31NO4/c1-3-5-6-7-8-9-10-14-22-16(19)15-12-11-13-18(15)17(20)21-4-2/l
InchiKey: RPQPMLBAGSABBH-UHFFFAOYSA-N
Formula: C17H31NO4
SMILES: CCCCCCCCCOC(=O)C1CCCN1C(=O)OCC
Mol. weight [g/mol]: 313.44

Physical Properties

Property code	Value	Unit	Source
hvap	97.97	kJ/mol	Cheméo Relay (1.0)
log10ws	-3.93		Cheméo Relay (1.0)
logp	3.901		Crippen Method
mcvol	264.390	ml/mol	McGowan Method
rinpol	2095.00		NIST Webbook
rinpol	2095.00		NIST Webbook
tb	615.04	K	Cheméo Relay (1.0)
tf	289.60	K	Cheméo Relay (1.0)

Sources

Cheméo Relay (1.0, beta):
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=U320841&Units=SI
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):

Legend

hvap: Enthalpy of vaporization at standard conditions
log10ws: Log10 of Water solubility in mol/l
logp: Octanol/Water partition coefficient

mcvol:	McGowan's characteristic volume
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tf:	Normal melting (fusion) point