

Padimate O

Other names:

2-Ethylhexyl 4-dimethylaminobenzoate
p-Dimethylaminobenzoic acid 2-ethylhexyl ester
Benzoic acid, 4-(dimethylamino)-, 2-ethylhexyl ester
2-Ethylhexyl p-(dimethylamino)benzoate
Arlatone UVB
Escalol 507
4-(Dimethylamino)benzoic acid 2-ethylhexyl ester
Eusolex 6007
Octyl dimethyl PABA

Inchi:

InChI=1S/C17H27NO2/c1-5-7-8-14(6-2)13-20-17(19)15-9-11-16(12-10-15)18(3)4/h9-12,17,20

InchiKey:

WYWZRNAHINYAEF-UHFFFAOYSA-N

Formula:

C17H27NO2

SMILES:

CCCCC(CC)COC(=O)c1ccc(N(C)C)cc1

Mol. weight [g/mol]:

277.40

CAS:

21245-02-3

Physical Properties

Property code	Value	Unit	Source
gf	69.46	kJ/mol	Joback Method
hf	-351.70	kJ/mol	Joback Method
hfus	35.72	kJ/mol	Joback Method
hvap	67.19	kJ/mol	Joback Method
log10ws	-4.31		Crippen Method
logp	4.126		Crippen Method
mcvol	244.050	ml/mol	McGowan Method
pc	1623.29	kPa	Joback Method
rinpol	2198.00		NIST Webbook
tb	708.31	K	Joback Method
tc	903.86	K	Joback Method
tf	409.92	K	Joback Method
vc	0.915	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	695.76	J/mol×K	708.31	Joback Method
cpg	713.53	J/mol×K	740.90	Joback Method
cpg	730.26	J/mol×K	773.49	Joback Method
cpg	745.99	J/mol×K	806.09	Joback Method
cpg	760.77	J/mol×K	838.68	Joback Method
cpg	774.61	J/mol×K	871.27	Joback Method
cpg	787.57	J/mol×K	903.86	Joback Method

Sources

Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C21245023&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.cheméo.com/doc/models/crippen_log10ws

Legend

cpg:	Ideal gas heat capacity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
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
pc:	Critical Pressure
rinpolar:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.cheméo.com/cid/44-034-5/Padimate-O.pdf>

Generated by Cheméo on 2024-04-26 04:44:17.236012305 +0000 UTC m=+16395906.156589620.

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