

Butyric acid, 4-phenyl-, heptyl ester

Inchi:	InChI=1S/C17H26O2/c1-2-3-4-5-9-15-19-17(18)14-10-13-16-11-7-6-8-12-16/h6-8,11-12H
InchiKey:	GZGITFJBUSHFNQ-UHFFFAOYSA-N
Formula:	C17H26O2
SMILES:	CCCCCCCOC(=O)CCc1ccccc1
Mol. weight [g/mol]:	262.39

Physical Properties

Property code	Value	Unit	Source
gf	-29.25	kJ/mol	Joback Method
hf	-402.48	kJ/mol	Joback Method
hfus	36.61	kJ/mol	Joback Method
hvap	64.87	kJ/mol	Joback Method
log10ws	-4.90		Crippen Method
logp	4.523		Crippen Method
mvol	234.070	ml/mol	McGowan Method
pc	1635.13	kPa	Joback Method
rinpol	1986.00		NIST Webbook
rinpol	1986.00		NIST Webbook
tb	691.33	K	Joback Method
tc	884.75	K	Joback Method
tf	379.93	K	Joback Method
vc	0.903	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	654.86	J/molxK	691.33	Joback Method
cpg	672.36	J/molxK	723.57	Joback Method
cpg	688.87	J/molxK	755.80	Joback Method
cpg	704.44	J/molxK	788.04	Joback Method
cpg	719.09	J/molxK	820.27	Joback Method
cpg	732.85	J/molxK	852.51	Joback Method
cpg	745.76	J/molxK	884.75	Joback Method
dvisc	0.0016807	Paxs	379.93	Joback Method

dvisc	0.0008086	Paxs	431.83	Joback Method
dvisc	0.0004552	Paxs	483.73	Joback Method
dvisc	0.0002864	Paxs	535.63	Joback Method
dvisc	0.0001956	Paxs	587.53	Joback Method
dvisc	0.0001421	Paxs	639.43	Joback Method
dvisc	0.0001083	Paxs	691.33	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=U406177&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

Legend

cp_g:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
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:	Log ₁₀ of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
m_cvol:	McGowan's characteristic volume
pc:	Critical Pressure
rinpol:	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.chemeo.com/cid/90-378-3/Butyric-acid-4-phenyl-heptyl-ester.pdf>

Generated by Cheméo on 2024-04-24 11:03:46.31368701 +0000 UTC m=+16245875.234264325.

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