

2,4-Bis(4-methoxyphenyl)-3,5-dimethyltetrahydrofuran

Inchi:	InChI=1S/C20H24O3/c1-13-19(15-5-9-17(21-3)10-6-15)14(2)23-20(13)16-7-11-18(22-4)
InchiKey:	MDWDSTVGZXXUID-UHFFFAOYSA-N
Formula:	C20H24O3
SMILES:	COc1ccc(C2OC(C)C(c3ccc(OC)cc3)C2C)cc1
Mol. weight [g/mol]:	312.40
CAS:	844679-34-1

Physical Properties

Property code	Value	Unit	Source
gf	40.38	kJ/mol	Joback Method
hf	-402.99	kJ/mol	Joback Method
hfus	42.36	kJ/mol	Joback Method
hvap	74.65	kJ/mol	Joback Method
log10ws	-5.08		Crippen Method
logp	4.583		Crippen Method
mvol	251.890	ml/mol	McGowan Method
pc	1641.76	kPa	Joback Method
rinpol	2502.80		NIST Webbook
tb	793.38	K	Joback Method
tc	1027.37	K	Joback Method
tf	462.25	K	Joback Method
vc	0.934	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	781.22	J/molxK	793.38	Joback Method
cpg	801.15	J/molxK	832.38	Joback Method
cpg	819.38	J/molxK	871.38	Joback Method
cpg	835.95	J/molxK	910.38	Joback Method
cpg	850.86	J/molxK	949.37	Joback Method
cpg	864.16	J/molxK	988.37	Joback Method
cpg	875.85	J/molxK	1027.37	Joback Method
dvisc	0.0007963	Paxs	462.25	Joback Method

dvisc	0.0005260	Paxs	517.44	Joback Method
dvisc	0.0003764	Paxs	572.63	Joback Method
dvisc	0.0002857	Paxs	627.82	Joback Method
dvisc	0.0002267	Paxs	683.00	Joback Method
dvisc	0.0001862	Paxs	738.19	Joback Method
dvisc	0.0001572	Paxs	793.38	Joback Method

Sources

Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
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=C844679341&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

cpg:	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:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	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/87-339-0/2-4-Bis-4-methoxyphenyl-3-5-dimethyltetrahydrofuran.pdf>

Generated by Cheméo on 2024-04-28 08:27:02.443188354 +0000 UTC m=+16582071.363765675.

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