

Other names

heptafluorobutylate 4,6-capable, 1,6-capable, 1,2,3,4,5,18,21,24,27,30,33,36,39,42-trideca-
 teradecaethylene glycol, bis(heptafluorobutylate)
 heptafluorobutylate

Property code	Value	Unit	Temperature [K]	Source
cpg	2280.65	J/mol×K	1437.52	Joback Method
cpg	2111.07	J/mol×K	1685.33	Joback Method
cpg	1812.16	J/mol×K	1933.14	Joback Method
cpg	1391.83	J/mol×K	2180.95	Joback Method
cpg	858.03	J/mol×K	2428.76	Joback Method
cpg	218.69	J/mol×K	2676.57	Joback Method
cpg	-518.24	J/mol×K	2924.38	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307I
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=U352022&Units=SI

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
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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