

2-[2-[2-[2-[2-(2,2,2-Trifluoroacetyl)oxyethoxy]et

Other names: Hexaethylene glycol bis(trifluoroacetate)
2,2,2-trifluoroacetate 20,20,20-Trifluoro-19-oxo-3,6,9,12,15,18-hexaoxonadec-1-yl trifluoroacetate

Inchi: InChI=1S/C16H24F6O9/c17-15(18,19)13(23)30-11-9-28-7-5-26-3-1-25-2-4-27-6-8-29-10

InchiKey: UUDKMFYUYZWJOD-UHFFFAOYSA-N

Formula: C16H24F6O9

SMILES: O=C(OCCOCCOCCOCCOCCOCCOC(=O)C(F)(F)F)C(F)(F)F

Mol. weight [g/mol]: 474.35

Physical Properties

Property code	Value	Unit	Source
gf	-2072.18	kJ/mol	Joback Method
hf	-2718.43	kJ/mol	Joback Method
hfus	52.36	kJ/mol	Joback Method
hvap	74.08	kJ/mol	Joback Method
log10ws	-1.01		Crippen Method
logp	1.280		Crippen Method
mcpvol	291.150	ml/mol	McGowan Method
pc	1126.83	kPa	Joback Method
rinpol	1941.70		NIST Webbook
rinpol	1941.70		NIST Webbook
tb	819.32	K	Joback Method
tc	1003.47	K	Joback Method
tf	533.93	K	Joback Method
vc	1.155	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	930.32	J/molxK	819.32	Joback Method
cpg	944.50	J/molxK	850.01	Joback Method
cpg	957.57	J/molxK	880.70	Joback Method
cpg	969.52	J/molxK	911.40	Joback Method
cpg	980.35	J/molxK	942.09	Joback Method
cpg	990.04	J/molxK	972.78	Joback Method

