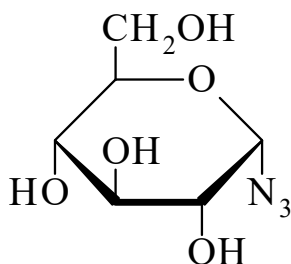




>>Entry Name: Glucopyranosyl azide, 9CI



α -D-Pyranose-form

CAS Registry Number: 122407-77-6

Molecular Formula: $C_6H_{11}N_3O_5$

Molecular Weight: 205.17

Accurate Mass: 205.069872

Percentage Composition: C 35.13%; H 5.40%; N 20.48%; O 38.99%

>>Variant: α -D-Pyranose-form

CAS Registry Number: 20379-60-6

Molecular Formula: $C_6H_{11}N_3O_5$

Molecular Weight: 205.17

Accurate Mass: 205.069872

Percentage Composition: C 35.13%; H 5.40%; N 20.48%; O 38.99%

Physical Description: Cryst. (MeCN/EtOH)

Melting Point: Mp 181.5-182.5° (176-178°)

Optical Rotation: $[\alpha]_D^{22}+258$ (c, 1.13 in H_2O)

>>Derivative: Tetra-Ac

Synonym(s): 2,3,4,6-Tetra-*O*-acetyl- α -D-glucopyranosyl azide

CAS Registry Number: 20379-61-7

Molecular Formula: $C_{14}H_{19}N_3O_9$

Molecular Weight: 373.319

Accurate Mass: 373.112132

Percentage Composition: C 45.04%; H 5.13%; N 11.26%; O 38.57%

Physical Description: Cryst. (EtOH)

Melting Point: Mp 100-101°

Optical Rotation: $[\alpha]_D^{21}+191.7$ (c, 0.532 in $CHCl_3$)

>>Derivative: 2,3,4-Tribenzyl

Synonym(s): 2,3,4-Tri-*O*-benzyl- α -D-glucopyranosylazide

CAS Registry Number: 87216-68-0

Molecular Formula: $C_{27}H_{29}N_3O_5$

Molecular Weight: 475.543

Accurate Mass: 475.210722

Percentage Composition: C 68.20%; H 6.15%; N 8.84%; O 16.82%

Melting Point: Mp 68-69°

Optical Rotation: $[\alpha]_D+96.87$ (c, 2.6 in $CHCl_3$)

>>Variant: β -D-Pyranose-form

CAS Registry Number: 20379-59-3

Molecular Formula: $C_6H_{11}N_3O_5$

Molecular Weight: 205.17

Accurate Mass: 205.069872

Percentage Composition: C 35.13%; H 5.40%; N 20.48%; O 38.99%

Physical Description: Cryst.

Melting Point: Mp 89° (86-88°)

Optical Rotation: $[\alpha]_D^{22}-28.5$ (c, 1.34 in H_2O)