



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:53 PM UTC

PDB ID : 1UG9 / pdb_00001ug9
Title : Crystal Structure of Glucodextranase from *Arthrobacter globiformis* I42
Authors : Mizuno, M.; Tonozuka, T.; Suzuki, S.; Uotsu-Tomita, R.; Ohtaki, A.; Kami-
tori, S.; Nishikawa, A.; Sakano, Y.
Deposited on : 2003-06-16
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

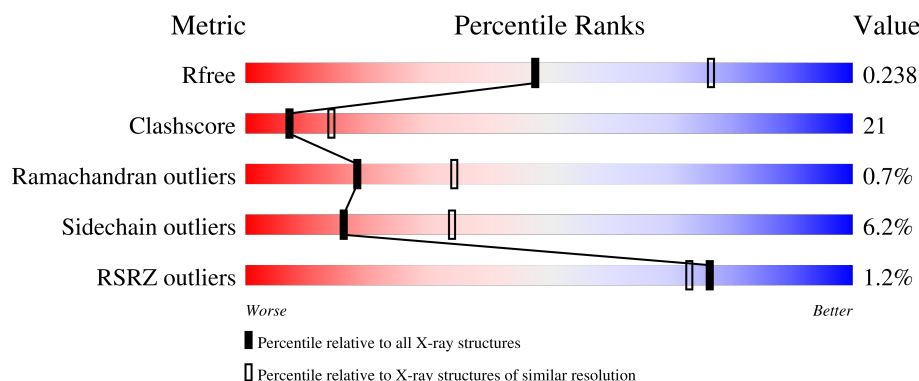
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	1020	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GOL	A	3001	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 8069 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called glucodextranase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1019	Total	C	N	O	S	0	0	0
			7506	4681	1286	1527	12			

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	6	Total	Ca	0	0
			6	6		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		

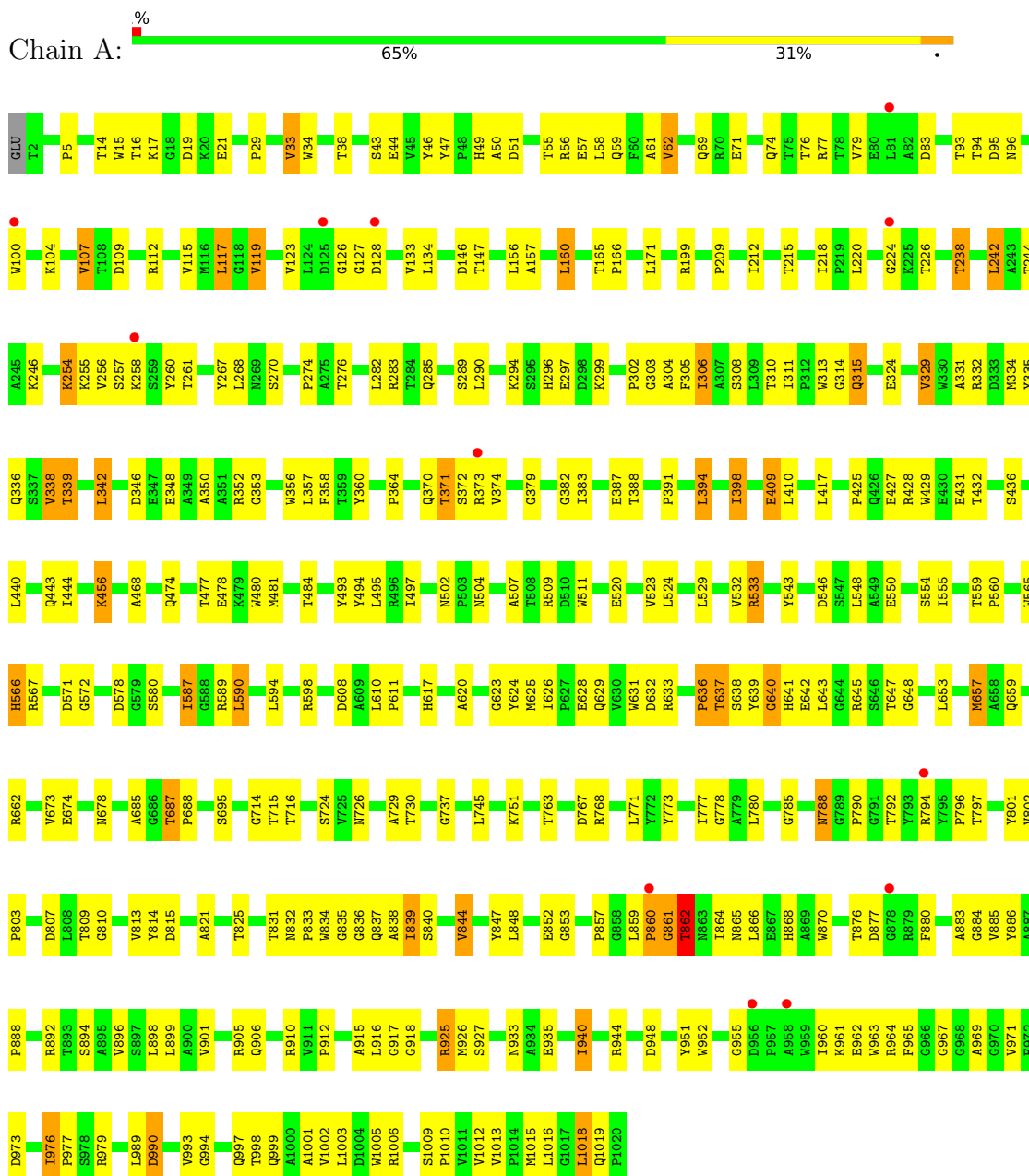
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	551	Total 551	O 551	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: glucodextranase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	199.86Å 88.53Å 80.78Å 90.00° 113.08° 90.00°	Depositor
Resolution (Å)	19.89 – 2.50 19.89 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.8 (19.89-2.50) 99.7 (19.89-2.50)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.53 (at 2.50Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.189 , 0.239 0.189 , 0.238	Depositor DCC
R_{free} test set	4522 reflections (10.08%)	wwPDB-VP
Wilson B-factor (Å ²)	22.6	Xtriage
Anisotropy	0.234	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 46.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.016 for -h-2*k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8069	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.26% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, CA

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.40	0/7683	0.91	17/10508 (0.2%)

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	126	GLY	N-CA-C	-7.01	105.68	114.16
1	A	687	THR	N-CA-C	6.92	118.43	109.64
1	A	360	TYR	N-CA-C	6.25	120.94	113.12
1	A	566	HIS	N-CA-C	-6.00	101.89	110.59
1	A	428	ARG	N-CA-C	5.82	119.48	112.38
1	A	388	THR	N-CA-C	-5.73	106.42	113.41
1	A	224	GLY	N-CA-C	5.72	119.14	112.50
1	A	119	VAL	N-CA-C	5.64	116.60	108.48
1	A	480	TRP	N-CA-C	5.59	118.57	111.69
1	A	329	VAL	N-CA-C	5.50	115.70	107.51
1	A	554	SER	N-CA-C	5.46	118.40	111.69
1	A	901	VAL	CA-C-N	5.41	125.48	119.32
1	A	901	VAL	C-N-CA	5.41	125.48	119.32
1	A	5	PRO	N-CA-C	-5.40	104.81	110.58
1	A	862	THR	N-CA-C	5.33	117.16	111.36
1	A	810	GLY	N-CA-C	5.32	118.25	110.38
1	A	647	THR	N-CA-C	-5.05	106.37	112.54

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	7506	0	7137	304	0
2	A	6	0	0	0	0
3	A	6	0	8	4	0
4	A	551	0	0	18	0
All	All	8069	0	7145	304	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 21.

All (304) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:831:THR:HG22	1:A:833:PRO:HD3	1.30	1.09
1:A:339:THR:HG21	1:A:662:ARG:HH21	1.31	0.94
1:A:71:GLU:HB2	1:A:93:THR:HG21	1.50	0.93
1:A:358:PHE:HB3	1:A:409:GLU:HG3	1.51	0.93
1:A:29:PRO:HA	1:A:299:LYS:HD2	1.52	0.92
1:A:104:LYS:HB3	1:A:117:LEU:HD21	1.52	0.91
1:A:285:GLN:HE21	1:A:617:HIS:HD2	1.14	0.91
1:A:993:VAL:HB	1:A:997:GLN:HG3	1.58	0.84
1:A:832:ASN:HD21	1:A:836:GLY:H	1.22	0.84
1:A:123:VAL:HG11	1:A:127:GLY:HA2	1.61	0.82
1:A:925:ARG:HB3	1:A:1015:MET:HG3	1.62	0.81
1:A:862:THR:CG2	1:A:864:ILE:HG12	2.10	0.81
1:A:342:LEU:HD21	1:A:398:ILE:CD1	2.11	0.81
1:A:17:LYS:HB3	1:A:38:THR:HG21	1.65	0.78
1:A:590:LEU:HD22	1:A:632:ASP:HB3	1.66	0.77
1:A:339:THR:HG23	1:A:662:ARG:HD3	1.64	0.77
1:A:865:ASN:ND2	1:A:1010:PRO:HG2	2.01	0.76
1:A:484:THR:H	1:A:504:ASN:HD21	1.32	0.76
1:A:315:GLN:H	1:A:315:GLN:NE2	1.83	0.76
1:A:324:GLU:OE2	1:A:371:THR:HG21	1.84	0.76
1:A:994:GLY:O	1:A:997:GLN:HG2	1.86	0.76
1:A:844:VAL:HG13	1:A:876:THR:HB	1.68	0.75
1:A:370:GLN:HE21	1:A:382:GLY:H	1.31	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:383:ILE:O	1:A:425:PRO:HG2	1.87	0.74
1:A:302:PRO:HB2	1:A:352:ARG:HD2	1.70	0.74
1:A:285:GLN:HE21	1:A:617:HIS:CD2	2.03	0.73
1:A:659:GLN:HE22	1:A:662:ARG:HH11	1.36	0.72
1:A:74:GLN:NE2	1:A:96:ASN:H	1.86	0.72
1:A:862:THR:HG21	1:A:864:ILE:HG12	1.72	0.71
1:A:255:LYS:HA	1:A:258:LYS:HE2	1.71	0.71
1:A:74:GLN:HE21	1:A:95:ASP:HA	1.55	0.70
1:A:507:ALA:O	1:A:520:GLU:HG3	1.92	0.70
1:A:726:ASN:HD21	1:A:751:LYS:H	1.40	0.69
1:A:342:LEU:HD11	1:A:398:ILE:HD11	1.75	0.69
1:A:861:GLY:HA3	1:A:940:ILE:O	1.93	0.69
1:A:305:PHE:HB2	1:A:334:MET:HG3	1.74	0.68
1:A:620:ALA:HB2	1:A:626:ILE:HD13	1.75	0.68
1:A:220:LEU:HD22	1:A:226:THR:HG21	1.77	0.67
1:A:383:ILE:HB	1:A:425:PRO:CG	2.25	0.67
1:A:456:LYS:NZ	1:A:456:LYS:HB2	2.10	0.66
1:A:777:ILE:HD11	1:A:815:ASP:HB2	1.77	0.66
1:A:313:TRP:HA	1:A:315:GLN:HE22	1.60	0.66
1:A:965:PHE:CE1	1:A:989:LEU:HD21	2.31	0.66
1:A:285:GLN:NE2	1:A:617:HIS:HD2	1.91	0.65
1:A:71:GLU:CB	1:A:93:THR:HG21	2.25	0.65
1:A:767:ASP:O	1:A:768:ARG:HD2	1.96	0.65
1:A:342:LEU:HD21	1:A:398:ILE:HD13	1.77	0.65
1:A:859:LEU:O	1:A:862:THR:HB	1.96	0.65
1:A:963:TRP:O	1:A:964:ARG:HD3	1.96	0.65
1:A:432:THR:HB	1:A:509:ARG:HH11	1.62	0.64
1:A:339:THR:CG2	1:A:662:ARG:HH21	2.07	0.64
1:A:14:THR:HG22	4:A:3169:HOH:O	1.97	0.64
1:A:50:ALA:HB1	1:A:306:ILE:HD12	1.78	0.64
1:A:567:ARG:HD3	1:A:571:ASP:OD2	1.99	0.63
1:A:566:HIS:HE1	1:A:632:ASP:OD1	1.81	0.63
1:A:529:LEU:HD12	3:A:3001:GOL:H11	1.80	0.63
1:A:339:THR:HG21	1:A:662:ARG:NH2	2.09	0.63
1:A:16:THR:HB	1:A:44:GLU:OE2	1.98	0.62
1:A:383:ILE:HB	1:A:425:PRO:HG2	1.82	0.62
1:A:831:THR:CG2	1:A:833:PRO:HD3	2.18	0.62
1:A:726:ASN:ND2	1:A:751:LYS:H	1.97	0.62
1:A:297:GLU:CD	1:A:352:ARG:HH21	2.07	0.61
1:A:342:LEU:HD21	1:A:398:ILE:HD11	1.83	0.60
1:A:289:SER:HA	1:A:657:MET:HG3	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:844:VAL:CG1	1:A:876:THR:HB	2.32	0.60
1:A:940:ILE:CG2	1:A:977:PRO:HB2	2.31	0.60
1:A:853:GLY:O	1:A:868:HIS:HD2	1.84	0.60
1:A:432:THR:HB	1:A:509:ARG:NH1	2.15	0.60
1:A:477:THR:HG22	1:A:494:TYR:OH	2.02	0.59
1:A:276:THR:HG23	4:A:3138:HOH:O	2.01	0.59
1:A:885:VAL:CG1	1:A:894:SER:HB3	2.32	0.59
1:A:594:LEU:HB3	3:A:3001:GOL:H2	1.84	0.59
1:A:336:GLN:O	1:A:339:THR:HG22	2.04	0.58
1:A:625:MET:HA	1:A:657:MET:HE1	1.86	0.58
1:A:999:GLN:HA	1:A:1002:VAL:CG1	2.34	0.58
1:A:884:GLY:HA2	1:A:896:VAL:HG13	1.84	0.58
1:A:334:MET:O	1:A:338:VAL:HG13	2.04	0.58
1:A:104:LYS:HB3	1:A:117:LEU:CD2	2.29	0.58
1:A:370:GLN:NE2	1:A:382:GLY:H	2.01	0.58
1:A:916:LEU:O	1:A:918:GLY:N	2.37	0.57
1:A:870:TRP:CH2	1:A:888:PRO:HA	2.38	0.57
1:A:631:TRP:HB3	1:A:642:GLU:O	2.04	0.57
1:A:1005:TRP:H	1:A:1005:TRP:CD1	2.21	0.57
1:A:802:VAL:HG13	1:A:803:PRO:HD2	1.86	0.57
1:A:885:VAL:HG12	1:A:894:SER:HB3	1.87	0.57
1:A:587:ILE:HG12	1:A:589:ARG:CZ	2.35	0.57
1:A:780:LEU:HG	1:A:1018:LEU:HD12	1.87	0.57
1:A:21:GLU:CD	1:A:79:VAL:HG21	2.30	0.56
1:A:160:LEU:HD12	1:A:166:PRO:HB2	1.87	0.56
1:A:998:THR:O	1:A:1002:VAL:HG12	2.05	0.56
1:A:999:GLN:HA	1:A:1002:VAL:HG12	1.88	0.56
1:A:315:GLN:H	1:A:315:GLN:CD	2.12	0.56
1:A:952:TRP:NE1	1:A:961:LYS:HA	2.21	0.56
1:A:839:ILE:HD13	1:A:839:ILE:H	1.71	0.55
1:A:555:ILE:C	1:A:555:ILE:HD12	2.31	0.55
1:A:628:GLU:HG2	1:A:629:GLN:HG3	1.88	0.55
1:A:976:ILE:HD13	1:A:976:ILE:H	1.71	0.55
1:A:282:LEU:HD12	4:A:3234:HOH:O	2.06	0.55
1:A:493:TYR:CD1	1:A:550:GLU:HG2	2.41	0.55
1:A:572:GLY:HA2	1:A:587:ILE:HA	1.89	0.55
1:A:371:THR:HG22	1:A:379:GLY:HA3	1.89	0.55
1:A:342:LEU:HD13	1:A:350:ALA:HB2	1.89	0.55
1:A:358:PHE:HB3	1:A:409:GLU:CG	2.32	0.55
1:A:62:VAL:HG11	1:A:100:TRP:CE2	2.42	0.54
1:A:156:LEU:HD13	1:A:238:THR:HG23	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:335:TYR:OH	1:A:533:ARG:NH1	2.37	0.54
1:A:852:GLU:O	1:A:868:HIS:HB3	2.07	0.54
1:A:165:THR:HG21	1:A:373:ARG:NH2	2.23	0.54
1:A:610:LEU:N	1:A:611:PRO:HD2	2.24	0.53
1:A:794:ARG:HD2	1:A:952:TRP:CZ3	2.43	0.53
1:A:297:GLU:HB2	1:A:305:PHE:CE2	2.43	0.53
1:A:57:GLU:CD	1:A:59:GLN:HE21	2.16	0.53
1:A:100:TRP:CD1	1:A:100:TRP:C	2.86	0.53
1:A:862:THR:HG22	1:A:864:ILE:H	1.73	0.53
1:A:51:ASP:OD2	1:A:374:VAL:HG22	2.09	0.53
1:A:133:VAL:HG23	1:A:218:ILE:CD1	2.39	0.53
1:A:788:ASN:C	1:A:788:ASN:HD22	2.17	0.53
1:A:859:LEU:HB3	1:A:860:PRO:HD2	1.90	0.53
1:A:578:ASP:OD2	1:A:580:SER:HB3	2.09	0.53
1:A:212:ILE:C	1:A:212:ILE:HD12	2.34	0.53
1:A:546:ASP:OD1	1:A:546:ASP:O	2.27	0.53
1:A:383:ILE:HB	1:A:425:PRO:HG3	1.90	0.53
1:A:905:ARG:HD3	4:A:3547:HOH:O	2.08	0.53
1:A:933:ASN:O	1:A:944:ARG:NH1	2.42	0.53
1:A:296:HIS:O	1:A:306:ILE:HD13	2.09	0.52
1:A:969:ALA:HB2	1:A:1005:TRP:CH2	2.44	0.52
1:A:578:ASP:O	1:A:637:THR:HB	2.09	0.52
1:A:348:GLU:CD	1:A:352:ARG:NH1	2.67	0.52
1:A:484:THR:HG22	1:A:484:THR:O	2.08	0.52
1:A:809:THR:OG1	1:A:825:THR:HG22	2.10	0.52
1:A:715:THR:HG23	1:A:737:GLY:O	2.10	0.52
1:A:254:LYS:O	1:A:258:LYS:HG3	2.09	0.52
1:A:146:ASP:O	1:A:209:PRO:HA	2.10	0.52
1:A:220:LEU:CD2	1:A:226:THR:HG21	2.40	0.51
1:A:555:ILE:HB	1:A:566:HIS:HB2	1.91	0.51
1:A:1006:ARG:HH21	1:A:1006:ARG:HG2	1.75	0.51
1:A:973:ASP:O	1:A:979:ARG:HA	2.09	0.51
1:A:16:THR:CG2	1:A:17:LYS:N	2.74	0.51
1:A:358:PHE:CB	1:A:409:GLU:HG3	2.33	0.51
1:A:74:GLN:HE21	1:A:96:ASN:H	1.55	0.51
1:A:50:ALA:CB	1:A:306:ILE:HD12	2.41	0.51
1:A:62:VAL:HG11	1:A:100:TRP:CZ2	2.45	0.51
1:A:832:ASN:ND2	1:A:840:SER:OG	2.42	0.51
1:A:289:SER:CA	1:A:657:MET:HG3	2.41	0.51
1:A:659:GLN:NE2	1:A:662:ARG:HH11	2.08	0.51
1:A:1015:MET:HE3	4:A:3255:HOH:O	2.09	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:631:TRP:NE1	1:A:633:ARG:HB2	2.25	0.50
1:A:778:GLY:HA3	1:A:1018:LEU:HB3	1.93	0.50
1:A:49:HIS:HB2	1:A:51:ASP:OD1	2.12	0.50
1:A:242:LEU:HD22	1:A:246:LYS:HD2	1.93	0.50
1:A:58:LEU:HA	1:A:134:LEU:O	2.12	0.49
1:A:659:GLN:HE22	1:A:662:ARG:NH1	2.07	0.49
1:A:71:GLU:HB2	1:A:93:THR:CG2	2.32	0.49
1:A:199:ARG:HH21	1:A:199:ARG:HG3	1.77	0.49
1:A:687:THR:HG23	1:A:688:PRO:HD2	1.94	0.49
1:A:695:SER:O	1:A:714:GLY:HA3	2.12	0.49
1:A:566:HIS:CE1	1:A:590:LEU:HD13	2.47	0.49
1:A:165:THR:HG21	1:A:373:ARG:HH21	1.77	0.49
1:A:484:THR:H	1:A:504:ASN:ND2	2.05	0.49
1:A:645:ARG:NH1	4:A:3023:HOH:O	2.46	0.49
1:A:857:PRO:HA	1:A:865:ASN:OD1	2.13	0.49
1:A:940:ILE:HG23	1:A:977:PRO:HB2	1.95	0.49
1:A:21:GLU:HB3	1:A:79:VAL:HG11	1.94	0.49
1:A:598:ARG:HB2	3:A:3001:GOL:H31	1.95	0.49
1:A:33:VAL:CG1	1:A:115:VAL:HG22	2.43	0.49
1:A:456:LYS:HB2	1:A:456:LYS:HZ2	1.77	0.48
1:A:866:LEU:HD12	1:A:927:SER:CB	2.43	0.48
1:A:357:LEU:HD13	1:A:391:PRO:HB3	1.95	0.48
1:A:864:ILE:HD12	1:A:1013:VAL:HG21	1.95	0.48
1:A:147:THR:O	1:A:157:ALA:HA	2.14	0.48
1:A:335:TYR:O	1:A:339:THR:HB	2.13	0.48
1:A:502:ASN:C	1:A:502:ASN:HD22	2.22	0.48
1:A:332:ARG:O	1:A:335:TYR:HB3	2.12	0.48
1:A:436:SER:HB2	1:A:497:ILE:HD12	1.94	0.48
1:A:745:LEU:HD12	1:A:745:LEU:N	2.28	0.48
1:A:785:GLY:HA2	4:A:3396:HOH:O	2.13	0.48
1:A:848:LEU:HD23	1:A:926:MET:HB3	1.96	0.48
1:A:364:PRO:HG2	4:A:3216:HOH:O	2.13	0.48
1:A:456:LYS:NZ	4:A:3523:HOH:O	2.35	0.47
1:A:456:LYS:HB2	1:A:456:LYS:HZ3	1.79	0.47
1:A:885:VAL:HG11	1:A:915:ALA:HB1	1.96	0.47
1:A:912:PRO:HG2	1:A:915:ALA:HB2	1.97	0.47
1:A:296:HIS:NE2	1:A:308:SER:HA	2.30	0.47
1:A:877:ASP:O	1:A:883:ALA:HB2	2.14	0.47
1:A:242:LEU:O	1:A:246:LYS:HG3	2.15	0.47
1:A:636:PRO:O	1:A:637:THR:O	2.33	0.47
1:A:16:THR:HG21	1:A:43:SER:HB3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:880:PHE:CZ	1:A:935:GLU:HG2	2.50	0.47
1:A:637:THR:O	1:A:638:SER:C	2.58	0.46
1:A:15:TRP:CG	1:A:311:ILE:HD13	2.51	0.46
1:A:55:THR:CG2	1:A:56:ARG:N	2.78	0.46
1:A:797:THR:OG1	1:A:962:GLU:HA	2.15	0.46
1:A:837:GLN:O	1:A:838:ALA:HB3	2.14	0.46
1:A:109:ASP:HA	1:A:256:VAL:HG13	1.96	0.46
1:A:989:LEU:O	1:A:990:ASP:HB3	2.16	0.46
1:A:294:LYS:HD3	1:A:294:LYS:O	2.15	0.46
1:A:1009:SER:HA	1:A:1010:PRO:C	2.40	0.46
1:A:831:THR:HG22	1:A:832:ASN:N	2.30	0.46
1:A:1019:GLN:NE2	4:A:3106:HOH:O	2.48	0.46
1:A:16:THR:O	1:A:314:GLY:HA3	2.16	0.46
1:A:33:VAL:HG11	1:A:115:VAL:HG22	1.96	0.46
1:A:862:THR:HG22	1:A:864:ILE:N	2.30	0.46
1:A:998:THR:HG23	1:A:1001:ALA:H	1.81	0.45
1:A:83:ASP:OD1	1:A:261:THR:HG21	2.16	0.45
1:A:468:ALA:HA	1:A:763:THR:OG1	2.17	0.45
1:A:796:PRO:HG2	1:A:801:TYR:CD1	2.51	0.45
1:A:802:VAL:HG11	4:A:3362:HOH:O	2.14	0.45
1:A:847:TYR:O	1:A:926:MET:HA	2.16	0.45
1:A:117:LEU:HD22	1:A:119:VAL:HG22	1.98	0.45
1:A:484:THR:N	1:A:504:ASN:HD21	2.06	0.45
1:A:270:SER:HA	4:A:3502:HOH:O	2.17	0.45
1:A:773:TYR:OH	1:A:906:GLN:NE2	2.50	0.45
1:A:796:PRO:HG2	1:A:801:TYR:CG	2.51	0.45
1:A:948:ASP:HB3	1:A:951:TYR:HB3	1.98	0.45
1:A:886:TYR:CE1	1:A:892:ARG:HB2	2.52	0.45
1:A:511:TRP:CE2	1:A:523:VAL:HG21	2.52	0.45
1:A:567:ARG:HD2	4:A:3069:HOH:O	2.17	0.45
1:A:636:PRO:C	1:A:637:THR:O	2.59	0.45
1:A:833:PRO:C	1:A:835:GLY:H	2.24	0.45
1:A:338:VAL:O	1:A:342:LEU:HB2	2.17	0.45
1:A:608:ASP:OD2	1:A:610:LEU:HD23	2.17	0.45
1:A:19:ASP:HB2	1:A:648:GLY:HA2	1.98	0.45
1:A:55:THR:HG22	1:A:56:ARG:N	2.33	0.44
1:A:940:ILE:HG21	1:A:977:PRO:HB2	1.99	0.44
1:A:440:LEU:HD23	1:A:481:MET:HE1	1.97	0.44
1:A:724:SER:HA	1:A:729:ALA:HA	1.99	0.44
1:A:34:TRP:HB2	1:A:46:TYR:HB2	2.00	0.44
1:A:802:VAL:CG1	1:A:803:PRO:HD2	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:107:VAL:HG21	1:A:257:SER:HA	1.99	0.44
1:A:127:GLY:O	1:A:128:ASP:HB2	2.17	0.44
1:A:860:PRO:C	1:A:862:THR:H	2.24	0.44
1:A:77:ARG:HG3	1:A:77:ARG:HH11	1.82	0.43
1:A:371:THR:HG23	1:A:372:SER:N	2.32	0.43
1:A:628:GLU:HB2	1:A:653:LEU:HD22	1.99	0.43
1:A:156:LEU:HD13	1:A:238:THR:CG2	2.48	0.43
1:A:267:TYR:OH	1:A:346:ASP:OD1	2.34	0.43
1:A:274:PRO:HG3	1:A:283:ARG:NE	2.33	0.43
1:A:440:LEU:HA	1:A:443:GLN:OE1	2.18	0.43
1:A:478:GLU:HG2	1:A:543:TYR:CD2	2.53	0.43
1:A:636:PRO:O	1:A:636:PRO:HG2	2.18	0.43
1:A:1012:VAL:O	1:A:1012:VAL:HG13	2.19	0.43
1:A:331:ALA:HB3	1:A:387:GLU:HB3	2.01	0.43
1:A:444:ILE:HD13	1:A:474:GLN:HA	2.00	0.43
1:A:315:GLN:H	1:A:315:GLN:HE21	1.60	0.43
1:A:61:ALA:HA	1:A:69:GLN:O	2.19	0.43
1:A:659:GLN:NE2	1:A:662:ARG:NH1	2.66	0.43
1:A:813:VAL:HG21	1:A:1018:LEU:HD21	1.99	0.43
1:A:590:LEU:HD22	1:A:632:ASP:CB	2.43	0.43
1:A:109:ASP:HA	1:A:256:VAL:CG1	2.49	0.43
1:A:133:VAL:HG23	1:A:218:ILE:HD13	1.99	0.43
1:A:559:THR:HG22	1:A:643:LEU:CD2	2.49	0.42
1:A:866:LEU:N	1:A:866:LEU:HD22	2.33	0.42
1:A:432:THR:HG23	4:A:3434:HOH:O	2.19	0.42
1:A:308:SER:OG	1:A:310:THR:HG22	2.19	0.42
1:A:348:GLU:HG2	1:A:352:ARG:NH1	2.35	0.42
1:A:598:ARG:HB2	3:A:3001:GOL:C3	2.49	0.42
1:A:715:THR:HG22	1:A:716:THR:N	2.35	0.42
1:A:792:THR:HG23	1:A:967:GLY:O	2.19	0.42
1:A:303:GLY:O	1:A:353:GLY:HA2	2.20	0.42
1:A:304:ALA:HB2	1:A:356:TRP:CE2	2.55	0.42
1:A:839:ILE:HD13	1:A:839:ILE:N	2.32	0.42
1:A:502:ASN:ND2	1:A:504:ASN:H	2.17	0.42
1:A:566:HIS:CE1	1:A:632:ASP:OD1	2.68	0.42
1:A:884:GLY:HA2	1:A:896:VAL:HG22	2.02	0.42
1:A:429:TRP:O	1:A:431:GLU:HG3	2.19	0.42
1:A:1006:ARG:HG2	1:A:1006:ARG:NH2	2.33	0.42
1:A:338:VAL:HG22	1:A:394:LEU:HD11	2.01	0.42
1:A:478:GLU:HG2	1:A:543:TYR:CG	2.54	0.42
1:A:623:GLY:O	1:A:624:TYR:HB2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:107:VAL:HG22	1:A:260:TYR:CG	2.55	0.41
1:A:559:THR:HB	1:A:560:PRO:HD2	2.02	0.41
1:A:639:TYR:O	1:A:640:GLY:C	2.61	0.41
1:A:899:LEU:HD22	1:A:910:ARG:HH22	1.85	0.41
1:A:47:TYR:CD1	1:A:47:TYR:C	2.98	0.41
1:A:502:ASN:C	1:A:502:ASN:ND2	2.77	0.41
1:A:631:TRP:HE1	1:A:633:ARG:HB2	1.84	0.41
1:A:673:VAL:O	1:A:673:VAL:HG22	2.19	0.41
1:A:790:PRO:HG2	1:A:792:THR:HG22	2.02	0.41
1:A:785:GLY:N	1:A:807:ASP:OD1	2.42	0.41
1:A:951:TYR:CE2	1:A:971:VAL:HA	2.55	0.41
1:A:560:PRO:HG2	4:A:3474:HOH:O	2.21	0.41
1:A:628:GLU:HB2	1:A:653:LEU:HD13	2.02	0.41
1:A:76:THR:HG23	1:A:94:THR:O	2.19	0.41
1:A:559:THR:HG22	1:A:643:LEU:HD21	2.01	0.41
1:A:112:ARG:HB2	1:A:244:THR:HG23	2.01	0.41
1:A:33:VAL:HG11	1:A:115:VAL:CG2	2.51	0.41
1:A:329:VAL:HG11	1:A:356:TRP:CH2	2.56	0.41
1:A:814:TYR:HB2	1:A:821:ALA:HB3	2.03	0.41
1:A:857:PRO:HG2	4:A:3326:HOH:O	2.20	0.41
1:A:960:ILE:HD13	1:A:960:ILE:HA	1.95	0.41
1:A:107:VAL:CG2	1:A:260:TYR:CD2	3.04	0.41
1:A:555:ILE:CD1	1:A:565:TRP:HD1	2.34	0.41
1:A:255:LYS:HB2	4:A:3039:HOH:O	2.21	0.40
1:A:631:TRP:CG	1:A:641:HIS:HB3	2.56	0.40
1:A:685:ALA:HB3	4:A:3083:HOH:O	2.20	0.40
1:A:477:THR:CG2	1:A:494:TYR:OH	2.68	0.40
1:A:484:THR:O	1:A:484:THR:CG2	2.69	0.40
1:A:860:PRO:O	1:A:862:THR:N	2.54	0.40
1:A:533:ARG:HH11	1:A:674:GLU:HG3	1.86	0.40
1:A:678:ASN:OD1	1:A:678:ASN:N	2.53	0.40
1:A:876:THR:HG23	1:A:898:LEU:HD22	2.03	0.40
1:A:976:ILE:HB	1:A:977:PRO:HD2	2.02	0.40
1:A:348:GLU:HG2	1:A:352:ARG:HH11	1.86	0.40
1:A:555:ILE:C	1:A:555:ILE:CD1	2.95	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	1017/1020 (100%)	966 (95%)	44 (4%)	7 (1%)	18	34

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	860	PRO
1	A	917	GLY
1	A	637	THR
1	A	861	GLY
1	A	636	PRO
1	A	640	GLY
1	A	955	GLY

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	752/753 (100%)	705 (94%)	47 (6%)	16	34

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	VAL
1	A	62	VAL
1	A	107	VAL
1	A	117	LEU

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Mol	Chain	Res	Type
1	A	160	LEU
1	A	171	LEU
1	A	215	THR
1	A	238	THR
1	A	242	LEU
1	A	254	LYS
1	A	268	LEU
1	A	290	LEU
1	A	306	ILE
1	A	315	GLN
1	A	338	VAL
1	A	339	THR
1	A	342	LEU
1	A	371	THR
1	A	394	LEU
1	A	398	ILE
1	A	409	GLU
1	A	410	LEU
1	A	417	LEU
1	A	427	GLU
1	A	456	LYS
1	A	495	LEU
1	A	524	LEU
1	A	532	VAL
1	A	533	ARG
1	A	548	LEU
1	A	587	ILE
1	A	590	LEU
1	A	657	MET
1	A	730	THR
1	A	771	LEU
1	A	788	ASN
1	A	834	TRP
1	A	839	ILE
1	A	844	VAL
1	A	862	THR
1	A	925	ARG
1	A	940	ILE
1	A	976	ILE
1	A	990	ASP
1	A	1003	LEU
1	A	1016	LEU

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Mol	Chain	Res	Type
1	A	1018	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	HIS
1	A	74	GLN
1	A	182	ASN
1	A	217	GLN
1	A	315	GLN
1	A	370	GLN
1	A	397	GLN
1	A	502	ASN
1	A	504	ASN
1	A	557	GLN
1	A	566	HIS
1	A	605	ASN
1	A	607	GLN
1	A	617	HIS
1	A	659	GLN
1	A	677	GLN
1	A	726	ASN
1	A	788	ASN
1	A	832	ASN
1	A	837	GLN
1	A	845	ASN
1	A	863	ASN
1	A	868	HIS
1	A	906	GLN
1	A	933	ASN
1	A	987	ASN
1	A	1019	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 6 are monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	GOL	A	3001	-	5,5,5	1.37	1 (20%)	5,5,5	0.64	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	A	3001	-	-	4/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	3001	GOL	C1-C2	2.15	1.60	1.51

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	3001	GOL	O1-C1-C2-C3
3	A	3001	GOL	O1-C1-C2-O2
3	A	3001	GOL	C1-C2-C3-O3
3	A	3001	GOL	O2-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	3001	GOL	4	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	1019/1020 (99%)	-0.27	12 (1%) 76 73	11, 25, 43, 67	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	224	GLY	4.2
1	A	125	ASP	2.9
1	A	956	ASP	2.5
1	A	100	TRP	2.4
1	A	81	LEU	2.2
1	A	860	PRO	2.1
1	A	958	ALA	2.1
1	A	128	ASP	2.1
1	A	373	ARG	2.1
1	A	794	ARG	2.0
1	A	258	LYS	2.0
1	A	878	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GOL	A	3001	6/6	0.68	0.33	61,62,63,64	0
2	CA	A	2002	1/1	0.96	0.04	33,33,33,33	0
2	CA	A	2006	1/1	0.97	0.04	32,32,32,32	0
2	CA	A	2004	1/1	0.98	0.04	36,36,36,36	0
2	CA	A	2005	1/1	0.99	0.02	28,28,28,28	0
2	CA	A	2003	1/1	0.99	0.03	30,30,30,30	0
2	CA	A	2001	1/1	0.99	0.03	28,28,28,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.