



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 04:22 AM UTC

PDB ID : 3WLR / pdb_00003wlr
Title : Crystal Structure Analysis of Plant Exohydrolase
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Deposited on : 2013-11-12
Resolution : 2.21 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

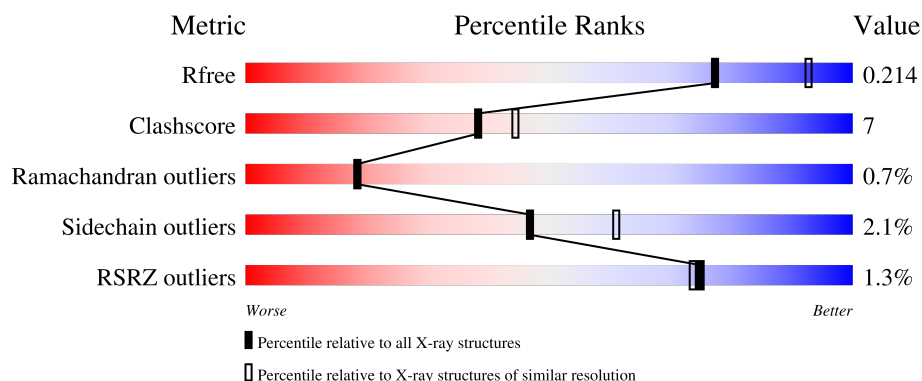
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.21 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	7682 (2.24-2.20)
Clashscore	190562	8402 (2.24-2.20)
Ramachandran outliers	187476	8303 (2.24-2.20)
Sidechain outliers	187428	8304 (2.24-2.20)
RSRZ outliers	180081	7683 (2.24-2.20)

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	609	 87% 12%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5273 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.

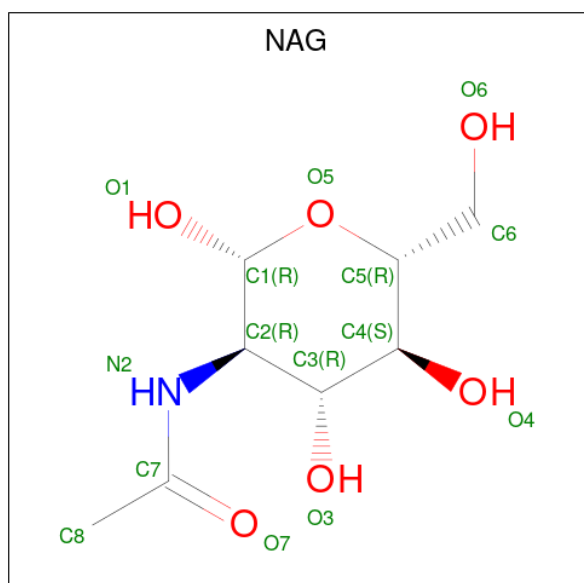
- Molecule 1 is a protein called Beta-D-glucan exohydrolase isoenzyme ExoI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	607	4621	2921	797	876	27	0	9	0

There are 7 discrepancies between the modelled and reference sequences:

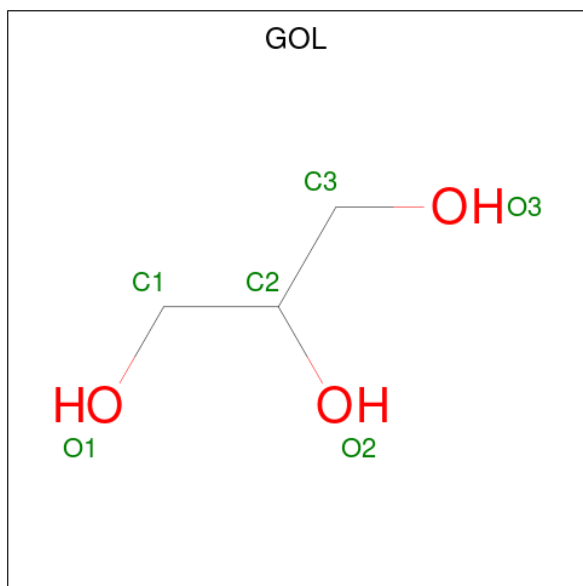
Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	HIS	-	expression tag	UNP Q9XEI3
A	-2	HIS	-	expression tag	UNP Q9XEI3
A	-1	ALA	-	expression tag	UNP Q9XEI3
A	0	ALA	-	expression tag	UNP Q9XEI3
A	158	ALA	ARG	engineered mutation	UNP Q9XEI3
A	161	ALA	GLU	engineered mutation	UNP Q9XEI3
A	320	LYS	ASN	SEE REMARK 999	UNP Q9XEI3

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

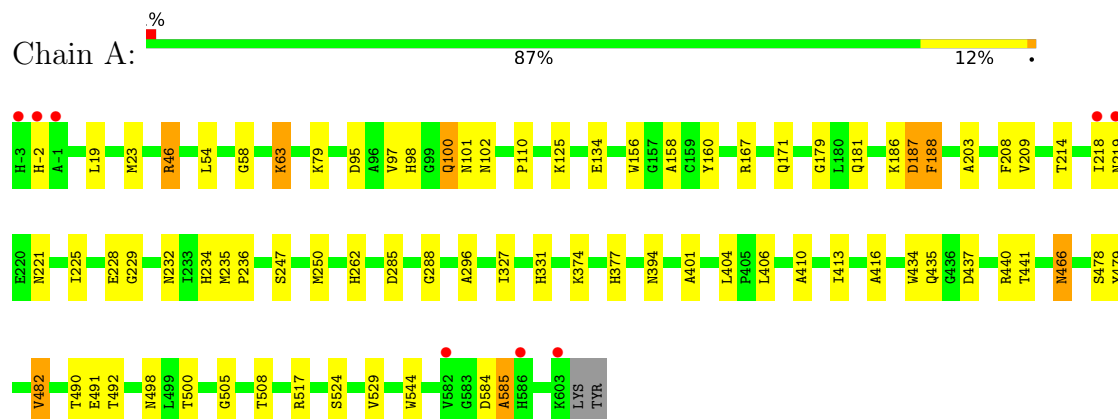
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	562	Total 562	O 562	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: Beta-D-glucan exohydrolase isoenzyme ExoI



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	100.69Å 100.69Å 183.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.37 – 2.21 46.37 – 2.21	Depositor EDS
% Data completeness (in resolution range)	98.8 (46.37-2.21) 99.2 (46.37-2.21)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.38 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, R_{free}	0.169 , 0.218 0.166 , 0.214	Depositor DCC
R_{free} test set	2405 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	49.0	Xtriage
Anisotropy	0.110	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 49.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5273	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.48% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	1.12	4/4765 (0.1%)	1.03	10/6474 (0.2%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	492	THR	CA-CB	5.92	1.62	1.53
1	A	98	HIS	CA-CB	5.47	1.57	1.53
1	A	179	GLY	N-CA	5.24	1.52	1.45
1	A	110	PRO	CA-C	5.16	1.58	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	188	PHE	N-CA-C	6.93	125.56	110.80
1	A	410	ALA	CA-C-N	-6.43	111.84	118.85
1	A	410	ALA	C-N-CA	-6.43	111.84	118.85
1	A	401	ALA	CA-C-N	-6.03	113.73	119.76
1	A	401	ALA	C-N-CA	-6.03	113.73	119.76
1	A	466[A]	ASN	CA-C-N	-5.92	113.86	120.14
1	A	466[A]	ASN	C-N-CA	-5.92	113.86	120.14
1	A	466[B]	ASN	CA-C-N	-5.92	113.86	120.14
1	A	466[B]	ASN	C-N-CA	-5.92	113.86	120.14
1	A	46	ARG	CB-CA-C	-5.49	102.44	110.95

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	4621	0	4593	59	0
2	A	42	0	39	1	0
3	A	48	0	64	11	0
4	A	562	0	0	15	0
All	All	5273	0	4696	65	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 7.

All (65) 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:46:ARG:HD3	4:A:809:HOH:O	1.33	1.26
1:A:394:ASN:HD21	1:A:404:LEU:H	1.14	0.94
1:A:491[B]:GLU:OE1	4:A:1299:HOH:O	1.84	0.94
1:A:466[A]:ASN:ND2	1:A:508:THR:OG1	2.01	0.94
1:A:498:ASN:HB3	4:A:1287:HOH:O	1.71	0.88
1:A:181:GLN:HE21	1:A:203:ALA:H	1.22	0.82
1:A:167:ARG:HH21	1:A:171:GLN:HE22	1.33	0.76
1:A:156:TRP:HE1	1:A:219:ASN:HD22	1.34	0.76
1:A:97:VAL:H	1:A:101:ASN:HD21	1.34	0.74
1:A:167:ARG:HH21	1:A:171:GLN:NE2	1.85	0.73
3:A:704:GOL:H31	4:A:1138:HOH:O	1.89	0.72
1:A:167:ARG:HH12	3:A:711:GOL:H31	1.55	0.71
1:A:100:GLN:HE21	1:A:100:GLN:HA	1.57	0.70
1:A:491[B]:GLU:CD	4:A:1299:HOH:O	2.35	0.64
1:A:95:ASP:OD1	3:A:705:GOL:H11	1.98	0.62
1:A:100:GLN:HE21	1:A:100:GLN:CA	2.15	0.60
3:A:704:GOL:H11	4:A:1321:HOH:O	2.02	0.60
1:A:58:GLY:H	1:A:102:ASN:ND2	2.00	0.59
1:A:181:GLN:HE22	1:A:247:SER:H	1.51	0.58
1:A:478:SER:O	1:A:517:ARG:NH1	2.37	0.58
1:A:491[B]:GLU:HG3	4:A:1269:HOH:O	2.03	0.58
1:A:517:ARG:HD3	4:A:1191:HOH:O	2.03	0.58
1:A:296:ALA:H	3:A:707:GOL:H2	1.69	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:181:GLN:NE2	1:A:203:ALA:H	1.98	0.56
1:A:327:ILE:HA	3:A:706:GOL:H31	1.89	0.55
1:A:134:GLU:OE2	1:A:377:HIS:HD2	1.90	0.55
1:A:187:ASP:OD2	1:A:187:ASP:N	2.39	0.53
1:A:524:SER:O	1:A:544:TRP:HA	2.09	0.53
1:A:225:ILE:HD11	1:A:229:GLY:HA3	1.91	0.53
1:A:434:TRP:O	1:A:490:THR:HG21	2.10	0.52
3:A:709:GOL:O3	3:A:709:GOL:O1	2.19	0.52
1:A:262:HIS:HE1	1:A:285:ASP:H	1.57	0.51
1:A:158:ALA:HB2	1:A:491[A]:GLU:HG3	1.93	0.51
1:A:167:ARG:NH2	1:A:171:GLN:HE22	2.04	0.51
1:A:331:HIS:HD2	4:A:801:HOH:O	1.92	0.51
1:A:97:VAL:H	1:A:101:ASN:ND2	2.03	0.50
3:A:709:GOL:H32	4:A:1113:HOH:O	2.11	0.49
1:A:262:HIS:CE1	1:A:288:GLY:HA3	2.46	0.49
1:A:435:GLN:HA	1:A:490:THR:HG23	1.95	0.48
1:A:214:THR:OG1	1:A:218:ILE:O	2.28	0.48
1:A:437:ASP:HB3	1:A:441:THR:HG21	1.94	0.48
1:A:500[A]:THR:HG22	1:A:529:VAL:HG21	1.96	0.47
1:A:250:MET:HE1	3:A:705:GOL:H12	1.95	0.47
1:A:134:GLU:OE2	1:A:377:HIS:CD2	2.69	0.46
1:A:416:ALA:O	1:A:482:VAL:HA	2.16	0.46
1:A:232[B]:ASN:CG	4:A:821:HOH:O	2.59	0.45
1:A:404:LEU:HD23	1:A:406:LEU:HG	1.98	0.45
1:A:100:GLN:O	1:A:100:GLN:HG3	2.17	0.45
1:A:156:TRP:HE1	1:A:219:ASN:ND2	2.09	0.44
1:A:125:LYS:NZ	4:A:984:HOH:O	2.49	0.44
3:A:707:GOL:H12	4:A:1052:HOH:O	2.18	0.44
1:A:235:MET:N	1:A:236:PRO:CD	2.81	0.43
1:A:221:ASN:HA	2:A:701:NAG:H83	1.99	0.43
1:A:500[A]:THR:HG22	1:A:529:VAL:CG2	2.49	0.43
1:A:63:LYS:HE3	4:A:859:HOH:O	2.19	0.43
1:A:19:LEU:O	1:A:23:MET:HG3	2.19	0.42
1:A:413:ILE:HG21	1:A:479:TYR:CZ	2.54	0.42
1:A:262:HIS:CE1	1:A:285:ASP:H	2.37	0.42
1:A:234:HIS:HE1	4:A:1115:HOH:O	2.04	0.41
1:A:100:GLN:HA	1:A:100:GLN:NE2	2.30	0.41
1:A:225:ILE:CD1	1:A:229:GLY:HA3	2.51	0.41
1:A:208:PHE:HA	1:A:209:VAL:HA	1.93	0.41
3:A:711:GOL:O3	3:A:711:GOL:O1	2.38	0.40
1:A:160:TYR:CD2	1:A:160:TYR:C	2.99	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:584:ASP:HB3	1:A:585:ALA:H	1.65	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	614/609 (101%)	585 (95%)	25 (4%)	4 (1%)	18 18

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	-2	HIS
1	A	188	PHE
1	A	585	ALA
1	A	505	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	495/488 (101%)	485 (98%)	10 (2%)	48 63

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	LEU
1	A	63	LYS
1	A	79	LYS
1	A	100	GLN
1	A	186	LYS
1	A	187	ASP
1	A	228	GLU
1	A	374	LYS
1	A	440	ARG
1	A	482	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	GLN
1	A	100	GLN
1	A	101	ASN
1	A	102	ASN
1	A	112	ASN
1	A	171	GLN
1	A	181	GLN
1	A	199	ASN
1	A	219	ASN
1	A	234	HIS
1	A	262	HIS
1	A	331	HIS
1	A	377	HIS
1	A	394	ASN
1	A	487	HIS

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

11 ligands are modelled in this entry.

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	710	-	5,5,5	0.40	0	5,5,5	1.25	1 (20%)
3	GOL	A	711	-	5,5,5	0.33	0	5,5,5	0.39	0
3	GOL	A	709	-	5,5,5	0.41	0	5,5,5	0.68	0
3	GOL	A	704	-	5,5,5	0.41	0	5,5,5	0.64	0
3	GOL	A	706	-	5,5,5	0.67	0	5,5,5	1.19	0
3	GOL	A	707	-	5,5,5	0.39	0	5,5,5	0.54	0
3	GOL	A	708	-	5,5,5	0.45	0	5,5,5	0.50	0
3	GOL	A	705	-	5,5,5	0.27	0	5,5,5	0.82	0
2	NAG	A	703	1	14,14,15	1.00	1 (7%)	17,19,21	2.69	8 (47%)
2	NAG	A	701	1	14,14,15	0.82	1 (7%)	17,19,21	1.01	1 (5%)
2	NAG	A	702	1	14,14,15	0.67	0	17,19,21	1.42	3 (17%)

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	710	-	-	4/4/4/4	-
3	GOL	A	711	-	-	2/4/4/4	-
3	GOL	A	709	-	-	4/4/4/4	-
3	GOL	A	704	-	-	2/4/4/4	-
3	GOL	A	706	-	-	2/4/4/4	-
3	GOL	A	707	-	-	0/4/4/4	-
3	GOL	A	708	-	-	4/4/4/4	-
3	GOL	A	705	-	-	2/4/4/4	-
2	NAG	A	703	1	-	3/6/23/26	0/1/1/1
2	NAG	A	701	1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	A	702	1	-	4/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	703	NAG	C2-N2	2.21	1.49	1.46
2	A	701	NAG	C1-C2	2.16	1.55	1.52

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	703	NAG	C2-N2-C7	8.23	133.93	122.90
2	A	702	NAG	C8-C7-N2	3.67	122.20	116.12
2	A	703	NAG	C1-O5-C5	2.81	115.95	112.19
2	A	703	NAG	C4-C3-C2	-2.77	106.96	111.02
2	A	703	NAG	O3-C3-C2	2.64	114.89	109.40
2	A	703	NAG	C1-C2-N2	2.47	114.32	110.43
2	A	701	NAG	O7-C7-C8	-2.35	117.88	122.05
2	A	703	NAG	C3-C4-C5	2.35	114.49	110.23
2	A	703	NAG	O5-C1-C2	-2.31	107.72	111.29
2	A	703	NAG	O7-C7-N2	2.22	125.90	121.98
2	A	702	NAG	C4-C3-C2	-2.11	107.92	111.02
2	A	702	NAG	O7-C7-N2	-2.07	118.33	121.98
3	A	710	GOL	O1-C1-C2	-2.01	101.33	110.38

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	703	NAG	C3-C2-N2-C7
3	A	704	GOL	C1-C2-C3-O3
3	A	709	GOL	O1-C1-C2-C3
3	A	709	GOL	C1-C2-C3-O3
3	A	709	GOL	O2-C2-C3-O3
3	A	710	GOL	O1-C1-C2-O2
3	A	710	GOL	O1-C1-C2-C3
2	A	702	NAG	O5-C5-C6-O6
2	A	702	NAG	C4-C5-C6-O6
2	A	701	NAG	C8-C7-N2-C2
2	A	701	NAG	O7-C7-N2-C2
2	A	702	NAG	C8-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
2	A	702	NAG	O7-C7-N2-C2
2	A	703	NAG	C4-C5-C6-O6
3	A	705	GOL	C1-C2-C3-O3
3	A	706	GOL	C1-C2-C3-O3
3	A	708	GOL	O1-C1-C2-C3
3	A	708	GOL	C1-C2-C3-O3
3	A	710	GOL	C1-C2-C3-O3
3	A	711	GOL	O1-C1-C2-C3
3	A	704	GOL	O2-C2-C3-O3
3	A	706	GOL	O2-C2-C3-O3
3	A	709	GOL	O1-C1-C2-O2
3	A	711	GOL	O1-C1-C2-O2
3	A	705	GOL	O1-C1-C2-O2
3	A	710	GOL	O2-C2-C3-O3
2	A	703	NAG	O5-C5-C6-O6
3	A	708	GOL	O2-C2-C3-O3
3	A	708	GOL	O1-C1-C2-O2

There are no ring outliers.

7 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	711	GOL	2	0
3	A	709	GOL	2	0
3	A	704	GOL	2	0
3	A	706	GOL	1	0
3	A	707	GOL	2	0
3	A	705	GOL	2	0
2	A	701	NAG	1	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	607/609 (99%)	-0.36	8 (1%) 75 73	33, 47, 65, 105	10 (1%)

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	-1	ALA	3.7
1	A	603	LYS	3.5
1	A	-3	HIS	3.3
1	A	582	VAL	2.9
1	A	-2	HIS	2.6
1	A	586	HIS	2.4
1	A	218	ILE	2.1
1	A	219	ASN	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
2	NAG	A	702	14/15	0.46	0.14	86,92,97,99	0
3	GOL	A	708	6/6	0.81	0.17	88,91,92,92	0
2	NAG	A	703	14/15	0.82	0.13	87,93,95,97	0
2	NAG	A	701	14/15	0.84	0.11	82,86,91,92	0
3	GOL	A	710	6/6	0.84	0.15	70,80,81,81	0
3	GOL	A	706	6/6	0.87	0.15	69,79,84,84	0
3	GOL	A	707	6/6	0.87	0.16	97,98,100,101	0
3	GOL	A	711	6/6	0.87	0.12	97,98,99,99	0
3	GOL	A	709	6/6	0.91	0.14	75,80,82,84	0
3	GOL	A	704	6/6	0.95	0.10	58,71,75,79	0
3	GOL	A	705	6/6	0.95	0.11	53,58,60,63	0

6.5 Other polymers [i](#)

There are no such residues in this entry.