



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 10:45 AM UTC

PDB ID : 4E8C / pdb_00004e8c
Title : Crystal structure of streptococcal beta-galactosidase in complex with galactose
Authors : Cheng, W.; Wang, L.; Bai, X.H.; Jiang, Y.L.; Li, Q.; Yu, G.; Zhou, C.Z.;
Chen, Y.X.
Deposited on : 2012-03-20
Resolution : 1.95 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

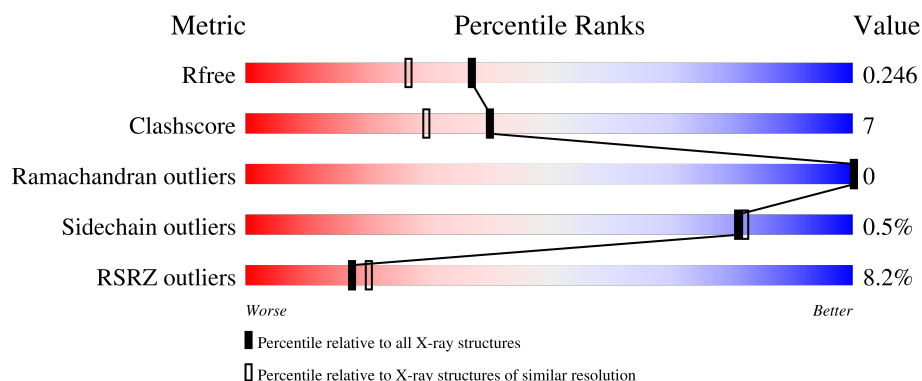
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	595	4% 86% 14%
1	B	595	12% 84% 15%

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	608	-	-	X	-

2 Entry composition [i](#)

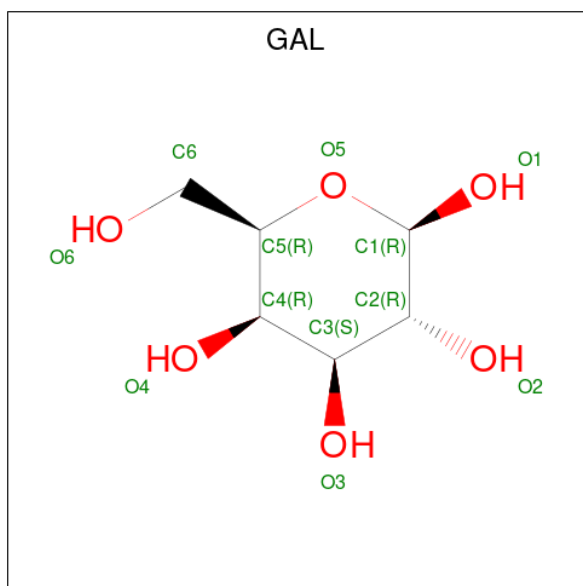
There are 4 unique types of molecules in this entry. The entry contains 10400 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Glycosyl hydrolase, family 35.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	594	Total	C	N	O	S	0	2	0
			4869	3140	799	907	23			
1	B	589	Total	C	N	O	S	0	0	0
			4825	3113	792	899	21			

- Molecule 2 is beta-D-galactopyranose (CCD ID: GAL) (formula: $C_6H_{12}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			12	6	6		
2	B	1	Total	C	O	0	0
			12	6	6		

- 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		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		

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

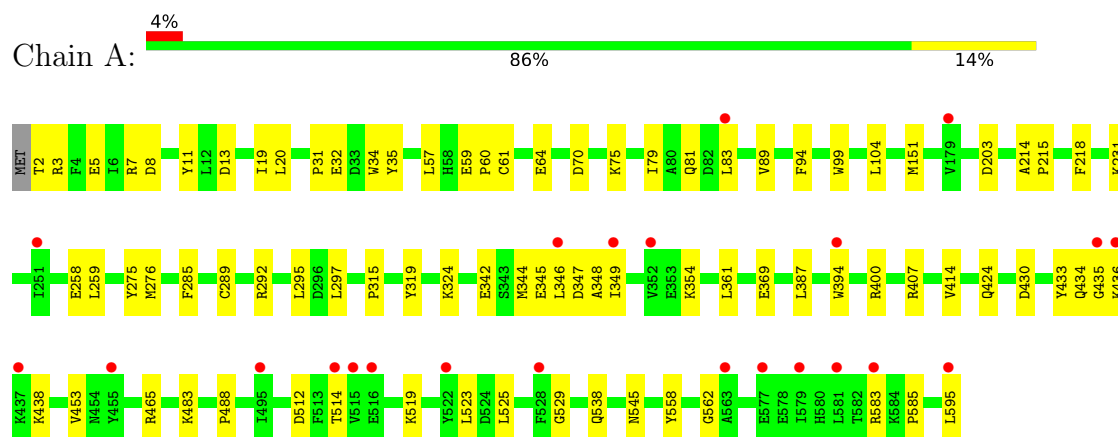
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	334	Total	O	0	0
			334	334		
4	B	258	Total	O	0	0
			258	258		

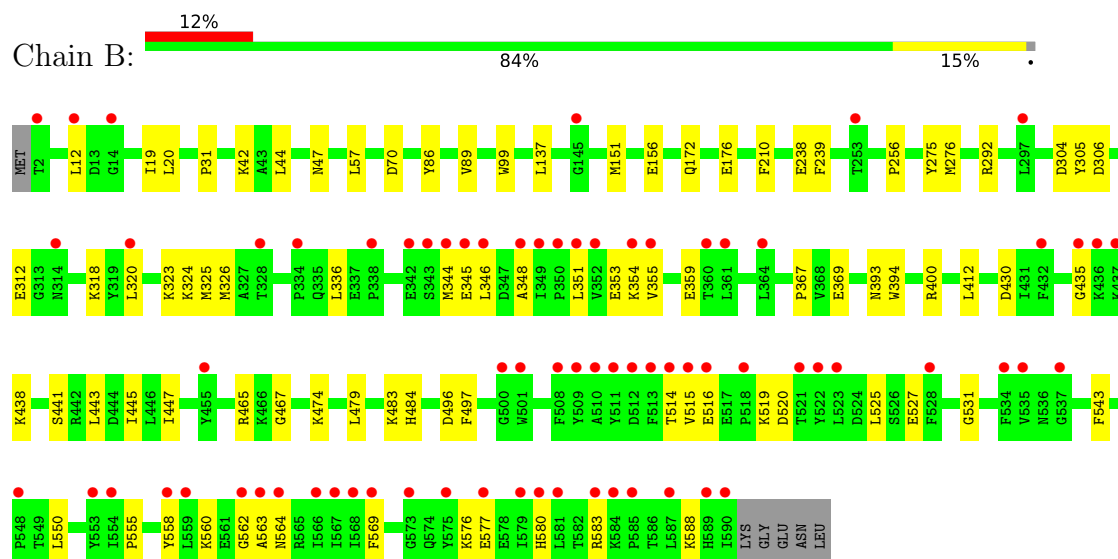
3 Residue-property plots [i](#)

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: Glycosyl hydrolase, family 35



- Molecule 1: Glycosyl hydrolase, family 35



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	80.35Å 82.37Å 99.60Å 90.00° 106.84° 90.00°	Depositor
Resolution (Å)	50.00 – 1.95 50.00 – 1.95	Depositor EDS
% Data completeness (in resolution range)	97.0 (50.00-1.95) 97.0 (50.00-1.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.44 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.5.0066	Depositor
R, R_{free}	0.205 , 0.243 0.216 , 0.246	Depositor DCC
R_{free} test set	4540 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	30.3	Xtriage
Anisotropy	0.387	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 44.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10400	wwPDB-VP
Average B, all atoms (Å ²)	40.0	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: GAL, GOL

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	0.42	0/5012	0.70	2/6789 (0.0%)
1	B	0.39	0/4962	0.72	1/6723 (0.0%)
All	All	0.40	0/9974	0.71	3/13512 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	516	GLU	N-CA-C	8.03	119.81	111.14
1	A	59	GLU	CA-C-N	5.23	125.03	119.28
1	A	59	GLU	C-N-CA	5.23	125.03	119.28

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	4869	0	4698	63	0
1	B	4825	0	4649	63	0
2	A	12	0	12	0	0
2	B	12	0	12	2	0
3	A	72	0	96	15	0
3	B	18	0	24	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	334	0	0	7	0
4	B	258	0	0	3	0
All	All	10400	0	9491	128	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 (128) 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:B:355:VAL:HG21	1:B:497:PHE:HA	1.27	1.15
1:B:355:VAL:CG2	1:B:497:PHE:HA	1.89	1.02
1:A:61[B]:CYS:SG	4:A:871:HOH:O	2.21	0.97
1:A:35:TYR:CE2	1:A:83:LEU:HD21	2.01	0.96
1:B:42:LYS:HE3	3:B:604:GOL:O3	1.71	0.91
1:A:538:GLN:OE1	3:A:608:GOL:H2	1.78	0.82
1:B:355:VAL:HG21	1:B:497:PHE:CA	2.10	0.80
1:A:558:TYR:CE2	3:A:608:GOL:H32	2.18	0.79
1:A:75:LYS:HE2	1:A:79:ILE:HD11	1.65	0.78
1:A:369:GLU:OE2	1:A:483:LYS:HE3	1.85	0.74
1:B:555:PRO:HG2	1:B:558:TYR:CD2	2.24	0.72
1:A:369:GLU:HG2	1:A:483:LYS:HG2	1.71	0.72
1:A:488:PRO:O	3:A:607:GOL:H12	1.90	0.71
1:B:42:LYS:CE	3:B:604:GOL:O3	2.43	0.67
1:A:292:ARG:NH1	1:A:297:LEU:HD11	2.10	0.67
1:A:324:LYS:NZ	4:A:925:HOH:O	2.28	0.66
1:A:354:LYS:O	3:A:609:GOL:H32	1.96	0.65
1:A:35:TYR:HE2	1:A:83:LEU:HD21	1.58	0.64
1:A:585:PRO:HG3	3:A:602:GOL:H31	1.81	0.62
1:A:345:GLU:HB3	1:A:583:ARG:HG3	1.83	0.61
1:B:465:ARG:HD2	4:B:930:HOH:O	2.01	0.60
1:A:5:GLU:HG3	1:A:11:TYR:HB2	1.82	0.60
1:B:344:MET:HE2	1:B:346:LEU:HD11	1.84	0.59
1:B:550:LEU:HD11	1:B:588:LYS:HB2	1.82	0.59
1:A:292:ARG:NH2	1:A:295:LEU:HD23	2.18	0.59
1:A:407:ARG:HH22	3:A:612:GOL:H31	1.67	0.58
1:B:172:GLN:O	1:B:176:GLU:HG2	2.03	0.58
1:B:345:GLU:C	1:B:346:LEU:HD12	2.28	0.58
1:A:5:GLU:CD	1:A:7:ARG:HH11	2.11	0.58
1:A:345:GLU:C	1:A:346:LEU:HD12	2.29	0.58
1:B:400:ARG:HD2	1:B:430:ASP:OD2	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:210:PHE:O	1:B:239:PHE:HA	2.04	0.57
1:B:355:VAL:HG23	1:B:497:PHE:HA	1.84	0.57
1:A:514:THR:HA	1:A:562:GLY:O	2.05	0.56
1:B:354:LYS:O	1:B:355:VAL:HG23	2.04	0.56
1:A:465:ARG:HD2	4:A:889:HOH:O	2.07	0.55
1:A:433:TYR:CE1	1:A:435:GLY:HA2	2.43	0.54
1:A:400:ARG:HD2	1:A:430:ASP:OD2	2.07	0.54
1:B:394:TRP:NE1	1:B:441:SER:OG	2.41	0.54
1:A:8:ASP:HA	1:A:231:LYS:O	2.09	0.53
1:A:285:PHE:HD2	3:A:607:GOL:H11	1.73	0.53
1:A:295:LEU:HD12	1:A:424:GLN:OE1	2.09	0.53
1:B:137:LEU:HD13	1:B:151:MET:HE1	1.90	0.53
3:A:613:GOL:H11	4:A:829:HOH:O	2.08	0.53
1:B:238:GLU:OE1	2:B:601:GAL:H1	2.09	0.53
1:A:558:TYR:CD2	3:A:608:GOL:H32	2.44	0.53
1:B:354:LYS:C	1:B:355:VAL:HG23	2.34	0.53
1:B:514:THR:HA	1:B:562:GLY:O	2.08	0.52
1:A:81:GLN:HG3	3:A:603:GOL:H12	1.91	0.52
1:B:354:LYS:O	1:B:355:VAL:CG2	2.57	0.52
1:B:351:LEU:HD23	1:B:577:GLU:OE2	2.09	0.52
1:A:595:LEU:C	1:A:595:LEU:HD23	2.34	0.52
1:A:394:TRP:CG	1:A:435:GLY:HA3	2.44	0.52
1:A:94:PHE:HB2	1:A:104:LEU:HD21	1.92	0.51
1:B:47:ASN:HB3	3:B:604:GOL:H12	1.92	0.50
1:B:344:MET:HB3	1:B:519:LYS:HB2	1.92	0.50
1:A:3:ARG:HD2	1:A:13:ASP:OD1	2.11	0.50
1:B:367:PRO:HA	1:B:484:HIS:O	2.12	0.50
1:A:529:GLY:HA2	1:A:545:ASN:ND2	2.27	0.50
1:B:137:LEU:HB3	1:B:151:MET:HE1	1.93	0.50
1:B:31:PRO:HG3	1:B:70:ASP:HA	1.94	0.49
1:B:355:VAL:CG2	1:B:497:PHE:CA	2.77	0.49
1:A:345:GLU:O	1:A:346:LEU:HD12	2.13	0.49
1:B:412:LEU:HD12	1:B:445:ILE:HG12	1.94	0.49
1:A:83:LEU:HG	4:A:912:HOH:O	2.13	0.49
1:A:394:TRP:CD1	1:A:435:GLY:HA3	2.49	0.48
1:B:564:ASN:HD22	1:B:564:ASN:N	2.11	0.48
1:B:325:MET:HE3	1:B:326:MET:N	2.28	0.48
1:B:527:GLU:OE1	1:B:576:LYS:NZ	2.43	0.48
1:B:292:ARG:NE	4:B:801:HOH:O	2.46	0.47
1:B:400:ARG:HD2	1:B:430:ASP:CG	2.38	0.47
1:A:259:LEU:C	1:A:259:LEU:HD13	2.38	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:GLN:HG3	3:A:603:GOL:C1	2.45	0.47
1:A:31:PRO:HA	1:A:34:TRP:CD2	2.50	0.47
1:B:323:LYS:NZ	1:B:336:LEU:O	2.42	0.47
1:A:214:ALA:HB3	1:A:215:PRO:HD3	1.97	0.46
1:B:44:LEU:HD12	1:B:323:LYS:HG2	1.97	0.46
1:A:31:PRO:HA	1:A:34:TRP:CE2	2.50	0.46
1:B:394:TRP:CD1	1:B:435:GLY:HA3	2.50	0.46
1:B:348:ALA:H	1:B:580:HIS:CE1	2.35	0.45
1:A:436:LYS:HG2	1:A:438:LYS:H	1.83	0.44
1:A:19:ILE:C	1:A:20:LEU:HD12	2.42	0.44
1:B:57:LEU:HD11	1:B:99:TRP:CZ2	2.52	0.44
1:B:474:LYS:HD3	1:B:479:LEU:HD21	1.98	0.44
1:B:515:VAL:HG22	1:B:564:ASN:ND2	2.33	0.44
1:A:344:MET:HE3	1:A:346:LEU:HD13	1.99	0.43
1:A:400:ARG:HD2	1:A:430:ASP:CG	2.42	0.43
1:B:156:GLU:CD	2:B:601:GAL:O1	2.61	0.43
1:B:137:LEU:HD13	1:B:151:MET:CE	2.48	0.43
1:B:256:PRO:HB3	1:B:318:LYS:HA	2.00	0.43
1:B:525:LEU:HD13	1:B:543:PHE:HB3	2.01	0.43
1:B:275:TYR:HA	1:B:276:MET:HA	1.82	0.43
1:B:563:ALA:C	1:B:564:ASN:HD22	2.27	0.43
1:B:305:TYR:O	1:B:306:ASP:HB3	2.19	0.43
1:B:12:LEU:HD21	1:B:86:TYR:OH	2.19	0.43
1:B:89:VAL:HB	1:B:151:MET:HE2	2.00	0.43
1:B:531:GLY:HA3	1:B:569:PHE:O	2.19	0.42
1:A:523:LEU:HD23	1:A:525:LEU:HD11	2.00	0.42
1:B:345:GLU:HB2	1:B:583:ARG:HG3	1.99	0.42
1:A:60:PRO:HD2	1:A:64:GLU:O	2.20	0.42
1:A:89:VAL:O	1:A:151:MET:HA	2.19	0.42
1:A:214:ALA:O	1:A:218:PHE:HB2	2.19	0.42
1:A:349:ILE:HG23	1:A:512:ASP:O	2.19	0.42
1:B:320:LEU:O	1:B:324:LYS:HG3	2.20	0.42
1:A:414:VAL:HG11	1:A:433:TYR:CE1	2.55	0.41
1:B:369:GLU:HG2	1:B:483:LYS:HG2	2.02	0.41
1:B:515:VAL:CG2	1:B:560:LYS:O	2.68	0.41
1:A:289:CYS:HB3	1:A:453:VAL:HA	2.02	0.41
1:A:342:GLU:H	1:A:342:GLU:HG2	1.65	0.41
1:B:447:ILE:HG21	1:B:467:GLY:HA2	2.03	0.41
1:A:258:GLU:HG2	3:A:610:GOL:H2	2.03	0.41
1:A:344:MET:HB3	1:A:519:LYS:HB2	2.01	0.41
1:A:32:GLU:OE1	3:A:608:GOL:O3	2.38	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:361:LEU:HD11	1:A:387:LEU:CD2	2.51	0.41
1:A:347:ASP:O	1:A:348:ALA:C	2.64	0.41
1:A:434:GLN:NE2	4:A:973:HOH:O	2.48	0.41
1:B:312:GLU:OE1	1:B:520:ASP:OD2	2.39	0.41
1:B:583:ARG:N	4:B:810:HOH:O	2.54	0.41
1:A:70:ASP:OD1	3:A:608:GOL:H12	2.21	0.41
1:B:355:VAL:HG21	1:B:496:ASP:O	2.21	0.40
1:B:359:GLU:OE1	1:B:359:GLU:HA	2.21	0.40
1:A:275:TYR:HA	1:A:276:MET:HA	1.83	0.40
1:B:19:ILE:C	1:B:20:LEU:HD12	2.46	0.40
1:A:315:PRO:HB3	1:A:319:TYR:CD2	2.57	0.40
3:A:613:GOL:H31	4:A:778:HOH:O	2.19	0.40
1:B:412:LEU:HD11	1:B:443:LEU:HD11	2.03	0.40
1:A:57:LEU:HD11	1:A:99:TRP:CZ2	2.56	0.40
1:B:304:ASP:HB2	3:B:603:GOL:H32	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	594/595 (100%)	574 (97%)	20 (3%)	0	100	100
1	B	587/595 (99%)	563 (96%)	24 (4%)	0	100	100
All	All	1181/1190 (99%)	1137 (96%)	44 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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	519/518 (100%)	517 (100%)	2 (0%)	84	85
1	B	513/518 (99%)	510 (99%)	3 (1%)	78	79
All	All	1032/1036 (100%)	1027 (100%)	5 (0%)	81	82

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	THR
1	A	203	ASP
1	B	353	GLU
1	B	393	ASN
1	B	438	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	374	GLN
1	A	381	GLN
1	A	594	ASN

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

17 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	$\# Z > 2$
2	GAL	A	601	-	12,12,12	0.62	0	17,17,17	0.92	0
2	GAL	B	601	-	12,12,12	0.63	0	17,17,17	0.93	1 (5%)
3	GOL	A	610	-	5,5,5	0.33	0	5,5,5	0.50	0
3	GOL	B	604	-	5,5,5	0.42	0	5,5,5	0.30	0
3	GOL	A	612	-	5,5,5	0.28	0	5,5,5	0.70	0
3	GOL	A	603	-	5,5,5	0.44	0	5,5,5	0.31	0
3	GOL	A	605	-	5,5,5	0.38	0	5,5,5	0.37	0
3	GOL	B	603	-	5,5,5	0.38	0	5,5,5	0.32	0
3	GOL	A	607	-	5,5,5	0.40	0	5,5,5	0.44	0
3	GOL	A	613	-	5,5,5	0.38	0	5,5,5	0.27	0
3	GOL	A	609	-	5,5,5	0.38	0	5,5,5	0.32	0
3	GOL	A	611	-	5,5,5	0.35	0	5,5,5	0.41	0
3	GOL	A	602	-	5,5,5	0.38	0	5,5,5	0.20	0
3	GOL	B	602	-	5,5,5	0.41	0	5,5,5	0.28	0
3	GOL	A	604	-	5,5,5	0.35	0	5,5,5	0.43	0
3	GOL	A	608	-	5,5,5	0.43	0	5,5,5	0.41	0
3	GOL	A	606	-	5,5,5	0.35	0	5,5,5	0.38	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
2	GAL	A	601	-	-	1/2/22/22	0/1/1/1
2	GAL	B	601	-	-	1/2/22/22	0/1/1/1
3	GOL	A	610	-	-	0/4/4/4	-
3	GOL	B	604	-	-	1/4/4/4	-
3	GOL	A	612	-	-	0/4/4/4	-
3	GOL	A	603	-	-	0/4/4/4	-
3	GOL	A	605	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	B	603	-	-	2/4/4/4	-
3	GOL	A	607	-	-	2/4/4/4	-
3	GOL	A	613	-	-	0/4/4/4	-
3	GOL	A	609	-	-	2/4/4/4	-
3	GOL	A	611	-	-	4/4/4/4	-
3	GOL	A	602	-	-	2/4/4/4	-
3	GOL	B	602	-	-	2/4/4/4	-
3	GOL	A	604	-	-	4/4/4/4	-
3	GOL	A	608	-	-	4/4/4/4	-
3	GOL	A	606	-	-	0/4/4/4	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	GAL	C1-O5-C5	-2.17	109.46	113.65

There are no chirality outliers.

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	602	GOL	O1-C1-C2-C3
3	A	607	GOL	C1-C2-C3-O3
3	A	608	GOL	C1-C2-C3-O3
3	A	609	GOL	O1-C1-C2-C3
3	A	611	GOL	O1-C1-C2-O2
3	B	602	GOL	O1-C1-C2-O2
3	B	602	GOL	O1-C1-C2-C3
3	B	603	GOL	O1-C1-C2-C3
3	A	602	GOL	O1-C1-C2-O2
3	A	605	GOL	O1-C1-C2-O2
3	A	607	GOL	O2-C2-C3-O3
3	A	604	GOL	O1-C1-C2-C3
3	A	605	GOL	O1-C1-C2-C3
3	A	608	GOL	O1-C1-C2-C3
3	A	611	GOL	O1-C1-C2-C3
3	A	611	GOL	C1-C2-C3-O3
3	A	608	GOL	O1-C1-C2-O2
3	A	608	GOL	O2-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
3	A	609	GOL	O1-C1-C2-O2
3	B	603	GOL	O1-C1-C2-O2
2	B	601	GAL	O5-C5-C6-O6
2	A	601	GAL	O5-C5-C6-O6
3	A	611	GOL	O2-C2-C3-O3
3	A	604	GOL	O1-C1-C2-O2
3	A	604	GOL	O2-C2-C3-O3
3	B	604	GOL	O1-C1-C2-O2
3	A	604	GOL	C1-C2-C3-O3

There are no ring outliers.

11 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	601	GAL	2	0
3	A	610	GOL	1	0
3	B	604	GOL	3	0
3	A	612	GOL	1	0
3	A	603	GOL	2	0
3	B	603	GOL	1	0
3	A	607	GOL	2	0
3	A	613	GOL	2	0
3	A	609	GOL	1	0
3	A	602	GOL	1	0
3	A	608	GOL	5	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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	594/595 (99%)	0.58	23 (3%)	43 50	23, 36, 52, 70	4 (0%)
1	B	589/595 (98%)	0.93	74 (12%)	8 9	27, 41, 59, 71	0
All	All	1183/1190 (99%)	0.75	97 (8%)	17 20	23, 38, 56, 71	4 (0%)

All (97) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	513	PHE	5.7
1	B	522	TYR	5.5
1	B	558	TYR	4.6
1	B	352	VAL	4.4
1	B	515	VAL	4.3
1	B	346	LEU	3.9
1	B	523	LEU	3.7
1	A	595	LEU	3.6
1	B	351	LEU	3.6
1	B	590	ILE	3.5
1	B	566	ILE	3.4
1	A	352	VAL	3.4
1	B	344	MET	3.4
1	B	2	THR	3.3
1	B	564	ASN	3.3
1	B	521	THR	3.2
1	A	455	TYR	3.2
1	B	350	PRO	3.2
1	B	320	LEU	3.1
1	A	346	LEU	3.1
1	B	569	PHE	3.1
1	B	514	THR	3.1
1	B	581	LEU	3.0
1	B	573	GLY	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	432	PHE	2.9
1	A	436	LYS	2.8
1	A	514	THR	2.8
1	A	515	VAL	2.8
1	B	585	PRO	2.8
1	B	580	HIS	2.8
1	B	516	GLU	2.8
1	B	575	TYR	2.7
1	B	253	THR	2.7
1	B	577	GLU	2.7
1	A	394	TRP	2.7
1	B	508	PHE	2.7
1	B	509	TYR	2.6
1	B	354	LYS	2.6
1	B	435	GLY	2.6
1	B	355	VAL	2.6
1	B	579	ILE	2.6
1	B	348	ALA	2.6
1	B	511	TYR	2.5
1	B	568	ILE	2.5
1	B	343	SER	2.5
1	B	455	TYR	2.5
1	B	537	GLY	2.5
1	B	589	HIS	2.5
1	B	563	ALA	2.5
1	B	587	LEU	2.4
1	A	516	GLU	2.4
1	B	554	ILE	2.4
1	B	328	THR	2.4
1	B	360	THR	2.4
1	B	510	ALA	2.4
1	B	535	VAL	2.4
1	A	83	LEU	2.4
1	B	553	TYR	2.4
1	B	559	LEU	2.3
1	B	534	PHE	2.3
1	B	562	GLY	2.3
1	B	361	LEU	2.3
1	A	528	PHE	2.3
1	A	583	ARG	2.3
1	B	548	PRO	2.3
1	A	251	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	495	ILE	2.2
1	B	436	LYS	2.2
1	B	528	PHE	2.2
1	B	437	LYS	2.2
1	A	179	VAL	2.2
1	A	435	GLY	2.2
1	B	500	GLY	2.2
1	B	518	PRO	2.2
1	A	563	ALA	2.2
1	A	579	ILE	2.2
1	B	584	LYS	2.2
1	A	581	LEU	2.2
1	B	297	LEU	2.2
1	B	338	PRO	2.2
1	B	342	GLU	2.1
1	B	567	ILE	2.1
1	A	437	LYS	2.1
1	B	583	ARG	2.1
1	B	345	GLU	2.1
1	B	349	ILE	2.1
1	B	512	ASP	2.1
1	B	364	LEU	2.1
1	A	349	ILE	2.1
1	A	577	GLU	2.1
1	B	334	PRO	2.1
1	A	522	TYR	2.0
1	B	14	GLY	2.0
1	B	145	GLY	2.0
1	B	12	LEU	2.0
1	B	501	TRP	2.0
1	B	314	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(Å ²)	Q<0.9
3	GOL	A	609	6/6	0.62	0.17	48,53,57,57	0
3	GOL	B	603	6/6	0.66	0.22	49,52,53,62	0
3	GOL	A	608	6/6	0.71	0.16	45,48,49,50	0
3	GOL	B	604	6/6	0.76	0.14	40,48,53,58	0
3	GOL	A	605	6/6	0.77	0.14	41,46,47,57	0
3	GOL	A	610	6/6	0.77	0.15	43,49,52,53	0
3	GOL	A	607	6/6	0.78	0.14	41,43,47,54	0
3	GOL	A	613	6/6	0.78	0.15	35,44,46,54	0
3	GOL	A	611	6/6	0.80	0.12	47,48,53,55	0
3	GOL	A	612	6/6	0.82	0.14	37,48,51,55	0
3	GOL	A	604	6/6	0.82	0.12	48,49,54,54	0
3	GOL	A	606	6/6	0.84	0.11	46,48,52,56	0
3	GOL	A	602	6/6	0.85	0.12	44,46,51,54	0
3	GOL	A	603	6/6	0.85	0.14	44,49,53,58	0
2	GAL	A	601	12/12	0.86	0.11	30,37,46,50	0
3	GOL	B	602	6/6	0.86	0.13	51,52,55,58	0
2	GAL	B	601	12/12	0.87	0.10	33,36,45,47	0

6.5 Other polymers [i](#)

There are no such residues in this entry.