



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:15 PM UTC

PDB ID : 5GNY / pdb_00005gny
Title : The structure of WT Bgl6
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Deposited on : 2016-07-25
Resolution : 2.10 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

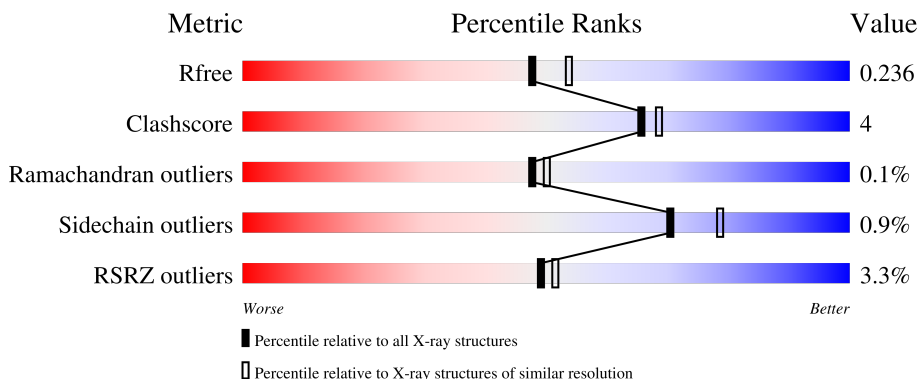
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	467	88% 8% .
1	B	467	10% 84% 11% .
1	C	467	88% 7% .
1	D	467	2% 84% 11% .

2 Entry composition [i](#)

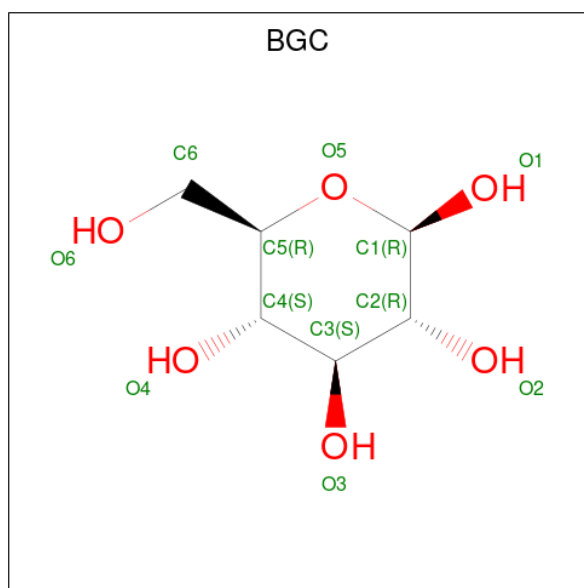
There are 3 unique types of molecules in this entry. The entry contains 15564 atoms, of which 0 are hydrogens and 0 are deuteriums.

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-glucosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	448	Total	C	N	O	S	0	0	0
			3580	2282	632	656	10			
1	B	448	Total	C	N	O	S	0	0	0
			3547	2263	629	645	10			
1	C	448	Total	C	N	O	S	0	0	0
			3582	2284	633	655	10			
1	D	448	Total	C	N	O	S	0	0	0
			3570	2275	630	655	10			

- Molecule 2 is beta-D-glucopyranose (CCD ID: BGC) (formula: C₆H₁₂O₆).



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		

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

- Molecule 3 is water.

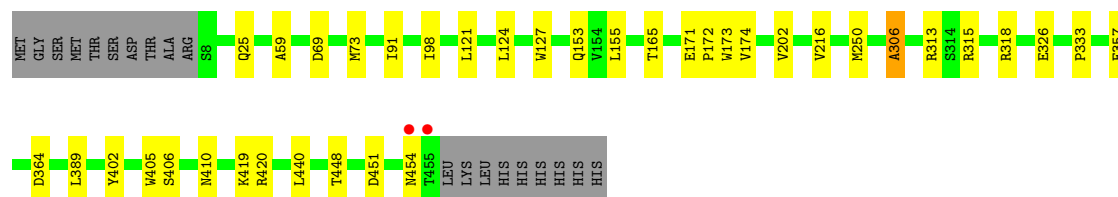
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	363	Total	O	0	0
			363	363		
3	B	211	Total	O	0	0
			211	211		
3	C	387	Total	O	0	0
			387	387		
3	D	276	Total	O	0	0
			276	276		

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.

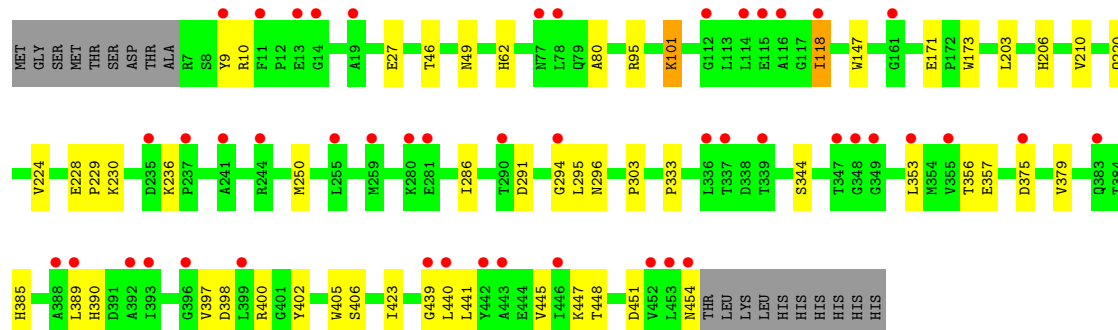
• Molecule 1: Beta-glucosidase

Chain A: 



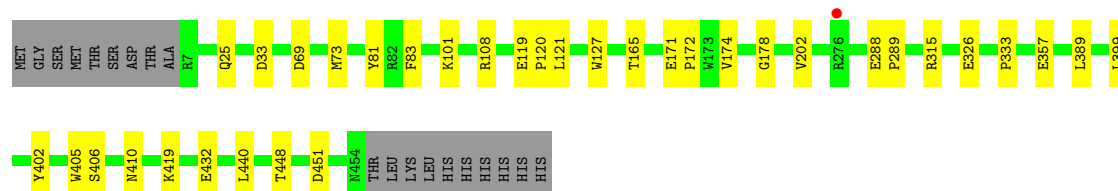
• Molecule 1: Beta-glucosidase

Chain B: 



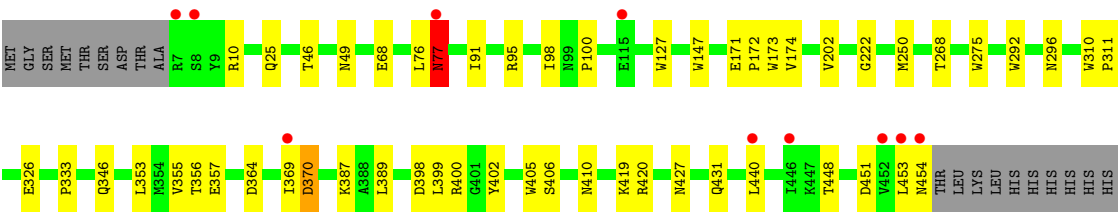
• Molecule 1: Beta-glucosidase

Chain C: 



• Molecule 1: Beta-glucosidase

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	68.36Å 159.70Å 197.36Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.97 – 2.10 41.97 – 2.10	Depositor EDS
% Data completeness (in resolution range)	98.7 (41.97-2.10) 89.3 (41.97-2.10)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.92 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.197 , 0.236 0.200 , 0.236	Depositor DCC
R_{free} test set	6250 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	22.0	Xtriage
Anisotropy	0.611	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 46.0	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	0.95	EDS
Total number of atoms	15564	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 41.68 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.2779e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: BGC

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	0.31	0/3692	0.71	4/5032 (0.1%)
1	B	0.31	0/3659	0.73	0/4993
1	C	0.31	0/3694	0.72	1/5033 (0.0%)
1	D	0.32	0/3682	0.74	3/5021 (0.1%)
All	All	0.31	0/14727	0.73	8/20079 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	D	0	1

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	333	PRO	N-CA-C	5.90	117.90	110.70
1	D	91	ILE	CA-C-N	5.50	125.59	119.32
1	D	91	ILE	C-N-CA	5.50	125.59	119.32
1	D	333	PRO	N-CA-C	5.28	117.14	110.70
1	A	333	PRO	N-CA-C	5.28	117.14	110.70
1	A	59	ALA	CB-CA-C	-5.18	110.58	116.54
1	A	306	ALA	CA-C-N	5.17	125.47	119.47
1	A	306	ALA	C-N-CA	5.17	125.47	119.47

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	369	ILE	Peptide

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	3580	0	3401	20	0
1	B	3547	0	3345	32	0
1	C	3582	0	3407	19	0
1	D	3570	0	3374	34	0
2	A	12	0	12	2	0
2	B	12	0	12	2	0
2	C	12	0	12	2	0
2	D	12	0	12	2	0
3	A	363	0	0	3	1
3	B	211	0	0	0	0
3	C	387	0	0	3	1
3	D	276	0	0	2	0
All	All	15564	0	13575	105	1

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 4.

All (105) 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:D:77:ASN:HD22	1:D:77:ASN:N	1.62	0.98
1:A:171:GLU:OE2	2:A:501:BGC:O1	1.97	0.81
1:D:448:THR:HB	1:D:451:ASP:HB2	1.73	0.69
1:A:448:THR:HB	1:A:451:ASP:HB2	1.74	0.69
1:C:171:GLU:OE2	2:C:501:BGC:O1	2.11	0.69
1:B:171:GLU:OE2	2:B:501:BGC:O1	2.13	0.66
1:D:171:GLU:OE2	2:D:501:BGC:H6C2	1.97	0.64
1:C:357:GLU:OE1	2:C:501:BGC:H1	2.00	0.62
1:B:10:ARG:O	1:B:390:HIS:NE2	2.33	0.61
1:D:76:LEU:C	1:D:77:ASN:HD22	2.10	0.60
1:B:296:ASN:OD1	1:B:356:THR:HG23	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:357:GLU:OE1	2:A:501:BGC:H1	2.04	0.58
1:B:398:ASP:OD2	1:B:400:ARG:NH2	2.35	0.58
1:B:9:TYR:N	1:B:454:ASN:OD1	2.37	0.57
1:D:296:ASN:OD1	1:D:356:THR:HG23	2.04	0.57
1:D:77:ASN:N	1:D:77:ASN:ND2	2.37	0.56
1:D:389:LEU:HD11	1:D:402:TYR:CD1	2.40	0.55
1:B:423:ILE:HB	1:B:439:GLY:CA	2.37	0.55
1:D:357:GLU:OE1	2:D:501:BGC:H4	2.08	0.53
1:C:448:THR:HB	1:C:451:ASP:HB2	1.91	0.52
1:A:318:ARG:NH1	3:A:610:HOH:O	2.42	0.52
1:C:172:PRO:HB3	1:C:202:VAL:HG11	1.91	0.52
1:D:173:TRP:CD1	1:D:250:MET:HA	2.45	0.52
1:C:389:LEU:HD11	1:C:402:TYR:CD1	2.46	0.51
1:A:91:ILE:HG12	1:A:98:ILE:HD13	1.91	0.51
1:D:398:ASP:OD2	1:D:400:ARG:NH2	2.43	0.51
1:C:121:LEU:HD22	1:C:165:THR:HB	1.93	0.51
1:D:364:ASP:HB2	1:D:420:ARG:HD2	1.93	0.51
1:B:357:GLU:OE1	2:B:501:BGC:H1	2.11	0.50
1:A:389:LEU:HD11	1:A:402:TYR:CD1	2.46	0.49
1:B:173:TRP:CD1	1:B:250:MET:HA	2.48	0.49
1:A:313:ARG:NH1	3:A:609:HOH:O	2.39	0.49
1:B:448:THR:HB	1:B:451:ASP:HB2	1.95	0.48
1:B:220:GLN:HB3	1:B:291:ASP:HB2	1.95	0.48
1:B:441:LEU:O	1:B:445:VAL:HG23	2.14	0.48
1:C:69:ASP:O	1:C:73:MET:HG3	2.14	0.48
1:B:398:ASP:CG	1:B:400:ARG:HH21	2.22	0.47
1:A:172:PRO:HB3	1:A:202:VAL:HG11	1.95	0.47
1:D:127:TRP:CH2	1:D:174:VAL:HG11	2.49	0.47
1:B:295:LEU:N	1:B:356:THR:HG22	2.29	0.47
1:D:268:THR:HG22	1:D:275:TRP:CD1	2.50	0.47
1:B:203:LEU:HB3	1:B:286:ILE:HG12	1.97	0.47
1:A:454:ASN:ND2	3:A:617:HOH:O	2.48	0.47
1:A:364:ASP:HB2	1:A:420:ARG:HD2	1.97	0.46
1:D:10:ARG:NH1	1:D:451:ASP:OD2	2.47	0.46
1:C:127:TRP:CH2	1:C:174:VAL:HG11	2.50	0.46
1:C:432:GLU:OE1	3:C:601:HOH:O	2.21	0.46
1:C:405:TRP:HA	1:C:406:SER:HA	1.61	0.46
1:D:222:GLY:HA3	1:D:292:TRP:O	2.16	0.46
1:A:153:GLN:HG3	1:A:216:VAL:HG21	1.98	0.46
1:A:25:GLN:O	1:A:410:ASN:HB2	2.16	0.46
1:A:326:GLU:O	1:A:419:LYS:HE3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:405:TRP:HA	1:A:406:SER:HA	1.59	0.45
1:D:405:TRP:HA	1:D:406:SER:HA	1.60	0.45
1:A:121:LEU:HD22	1:A:165:THR:HB	1.98	0.45
1:A:124:LEU:HD22	1:A:155:LEU:HD12	1.99	0.45
1:B:344:SER:HB2	1:B:397:VAL:HG13	1.99	0.45
1:C:178:GLY:HA2	3:C:898:HOH:O	2.17	0.45
1:D:46:THR:HB	1:D:49:ASN:ND2	2.33	0.44
1:D:25:GLN:O	1:D:410:ASN:HB2	2.18	0.44
1:C:81:TYR:CE2	1:C:83:PHE:HB3	2.53	0.43
1:D:172:PRO:HB3	1:D:202:VAL:HG11	2.00	0.43
1:A:306:ALA:HB2	1:A:315:ARG:HG2	1.99	0.43
1:D:451:ASP:C	1:D:454:ASN:HD22	2.26	0.43
1:A:69:ASP:O	1:A:73:MET:HG3	2.17	0.43
1:A:173:TRP:CD1	1:A:250:MET:HA	2.53	0.43
1:B:405:TRP:HA	1:B:406:SER:HA	1.57	0.43
1:D:451:ASP:HA	1:D:454:ASN:ND2	2.33	0.43
1:B:95:ARG:HD3	1:B:147:TRP:CD1	2.53	0.43
1:B:333:PRO:HG3	1:B:385:HIS:CE1	2.54	0.43
1:D:98:ILE:O	1:D:100:PRO:HD3	2.19	0.43
1:B:389:LEU:HD11	1:B:402:TYR:CD1	2.53	0.42
1:D:405:TRP:CD2	1:D:406:SER:HB2	2.54	0.42
1:B:224:VAL:HG22	1:B:294:GLY:HA3	2.01	0.42
1:D:68:GLU:OE1	3:D:601:HOH:O	2.21	0.42
1:B:356:THR:O	1:B:357:GLU:HG3	2.19	0.42
1:C:25:GLN:O	1:C:410:ASN:HB2	2.19	0.42
1:C:33:ASP:HB3	1:C:101:LYS:HD3	2.01	0.42
1:C:288:GLU:HA	1:C:289:PRO:HD3	1.90	0.42
1:B:206:HIS:O	1:B:210:VAL:HG23	2.19	0.42
1:C:389:LEU:HD22	1:C:399:LEU:HD21	2.01	0.42
1:A:127:TRP:CH2	1:A:174:VAL:HG11	2.55	0.42
1:B:10:ARG:NE	1:B:451:ASP:OD2	2.46	0.42
1:B:27:GLU:HA	1:B:62:HIS:HB3	2.01	0.42
1:B:101:LYS:HB2	1:B:101:LYS:HE2	1.72	0.42
1:B:228:GLU:HA	1:B:229:PRO:HD2	1.88	0.41
1:D:355:VAL:H	1:D:399:LEU:HD11	1.85	0.41
1:B:46:THR:HB	1:B:49:ASN:ND2	2.35	0.41
1:B:375:ASP:HB3	1:B:379:VAL:HG23	2.03	0.41
1:C:326:GLU:O	1:C:419:LYS:HE3	2.20	0.41
1:D:95:ARG:HD3	1:D:147:TRP:CD1	2.55	0.41
1:D:310:TRP:HA	1:D:311:PRO:HA	1.81	0.41
1:D:387:LYS:HE2	1:D:453:LEU:HB3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:108:ARG:NH2	3:C:604:HOH:O	2.29	0.41
1:D:427:ASN:O	1:D:431:GLN:N	2.50	0.41
1:B:296:ASN:OD1	1:B:357:GLU:HB2	2.21	0.41
1:B:296:ASN:CG	1:B:357:GLU:HB2	2.46	0.41
1:C:119:GLU:HA	1:C:120:PRO:HD3	1.93	0.41
1:D:326:GLU:O	1:D:419:LYS:HE3	2.21	0.41
1:B:80:ALA:C	1:B:118:ILE:HD11	2.45	0.40
1:B:230:LYS:HD2	1:B:303:PRO:HG3	2.03	0.40
1:D:405:TRP:CE2	1:D:406:SER:HB2	2.56	0.40
1:D:346:GLN:NE2	3:D:622:HOH:O	2.50	0.40
1:D:387:LYS:HG3	1:D:453:LEU:HD13	2.03	0.40
1:D:127:TRP:N	1:D:127:TRP:CD1	2.89	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:616:HOH:O	3:C:922:HOH:O[3_544]	2.16	0.04

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	446/467 (96%)	435 (98%)	11 (2%)	0	100	100
1	B	446/467 (96%)	434 (97%)	12 (3%)	0	100	100
1	C	446/467 (96%)	438 (98%)	8 (2%)	0	100	100
1	D	446/467 (96%)	433 (97%)	11 (2%)	2 (0%)	30	28
All	All	1784/1868 (96%)	1740 (98%)	42 (2%)	2 (0%)	48	50

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	370	ASP
1	D	77	ASN

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	360/378 (95%)	359 (100%)	1 (0%)	86	91
1	B	351/378 (93%)	345 (98%)	6 (2%)	53	62
1	C	360/378 (95%)	358 (99%)	2 (1%)	78	86
1	D	357/378 (94%)	353 (99%)	4 (1%)	65	74
All	All	1428/1512 (94%)	1415 (99%)	13 (1%)	70	78

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	440	LEU
1	B	101	LYS
1	B	118	ILE
1	B	236	LYS
1	B	353	LEU
1	B	440	LEU
1	B	447	LYS
1	C	315	ARG
1	C	440	LEU
1	D	77	ASN
1	D	353	LEU
1	D	370	ASP
1	D	440	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	358	ASN
1	A	380	HIS

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Mol	Chain	Res	Type
1	A	454	ASN
1	B	358	ASN
1	C	64	ASN
1	C	358	ASN
1	D	77	ASN
1	D	93	GLN
1	D	346	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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](#)

4 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	BGC	D	501	-	12,12,12	0.50	0	17,17,17	1.04	0
2	BGC	C	501	-	12,12,12	0.38	0	17,17,17	1.33	1 (5%)
2	BGC	B	501	-	12,12,12	0.47	0	17,17,17	1.18	1 (5%)
2	BGC	A	501	-	12,12,12	0.58	0	17,17,17	1.55	2 (11%)

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	BGC	D	501	-	-	2/2/22/22	0/1/1/1
2	BGC	C	501	-	-	0/2/22/22	0/1/1/1
2	BGC	B	501	-	-	2/2/22/22	0/1/1/1
2	BGC	A	501	-	-	1/2/22/22	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	BGC	C1-O5-C5	-4.79	104.38	113.65
2	C	501	BGC	C1-O5-C5	-3.68	106.54	113.65
2	B	501	BGC	C1-O5-C5	-3.31	107.25	113.65
2	A	501	BGC	O1-C1-C2	2.72	116.88	108.98

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	501	BGC	C4-C5-C6-O6
2	D	501	BGC	O5-C5-C6-O6
2	A	501	BGC	O5-C5-C6-O6
2	B	501	BGC	O5-C5-C6-O6
2	B	501	BGC	C4-C5-C6-O6

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	501	BGC	2	0
2	C	501	BGC	2	0
2	B	501	BGC	2	0
2	A	501	BGC	2	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	448/467 (95%)	-0.25	2 (0%) 88 90	14, 22, 34, 57	0
1	B	448/467 (95%)	0.92	47 (10%) 11 12	21, 39, 55, 62	0
1	C	448/467 (95%)	-0.23	1 (0%) 91 92	15, 22, 35, 49	0
1	D	448/467 (95%)	0.28	10 (2%) 62 65	15, 30, 45, 64	0
All	All	1792/1868 (95%)	0.18	60 (3%) 49 51	14, 26, 49, 64	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	439	GLY	4.5
1	B	440	LEU	4.2
1	D	77	ASN	4.0
1	B	19	ALA	3.7
1	B	453	LEU	3.6
1	B	118	ILE	3.6
1	B	11	PHE	3.4
1	C	276	ARG	3.3
1	B	454	ASN	3.1
1	B	383	GLN	3.1
1	B	347	THR	3.0
1	B	235	ASP	3.0
1	D	7	ARG	2.9
1	B	392	ALA	2.9
1	B	375	ASP	2.8
1	D	454	ASN	2.7
1	B	255	LEU	2.7
1	A	454	ASN	2.7
1	B	14	GLY	2.6
1	B	348	GLY	2.6
1	B	290	THR	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	259	MET	2.6
1	B	114	LEU	2.6
1	B	355	VAL	2.5
1	B	452	VAL	2.5
1	D	452	VAL	2.5
1	B	112	GLY	2.5
1	B	337	THR	2.5
1	B	349	GLY	2.5
1	B	396	GLY	2.5
1	B	393	ILE	2.4
1	B	446	ILE	2.4
1	B	353	LEU	2.4
1	B	237	PRO	2.4
1	B	115	GLU	2.3
1	B	281	GLU	2.3
1	B	389	LEU	2.3
1	D	453	LEU	2.3
1	B	388	ALA	2.3
1	B	161	GLY	2.3
1	B	280	LYS	2.3
1	A	455	THR	2.2
1	B	241	ALA	2.2
1	B	443	ALA	2.2
1	B	78	LEU	2.2
1	D	440	LEU	2.2
1	B	13	GLU	2.2
1	B	442	TYR	2.2
1	B	77	ASN	2.2
1	B	294	GLY	2.2
1	B	399	LEU	2.2
1	B	9	TYR	2.2
1	B	336	LEU	2.1
1	D	369	ILE	2.1
1	B	116	ALA	2.1
1	B	339	THR	2.0
1	D	446	ILE	2.0
1	D	115	GLU	2.0
1	D	8	SER	2.0
1	B	244	ARG	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	BGC	D	501	12/12	0.80	0.20	30,43,49,50	8
2	BGC	C	501	12/12	0.83	0.17	26,34,40,42	10
2	BGC	B	501	12/12	0.88	0.15	40,47,55,59	9
2	BGC	A	501	12/12	0.89	0.17	27,34,45,49	9

6.5 Other polymers [i](#)

There are no such residues in this entry.