



Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 10:17 pm BST

PDB ID : 6D1P
Title : Apo structure of Bacteroides uniformis beta-glucuronidase 3
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Deposited on : 2018-04-12
Resolution : 2.35 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

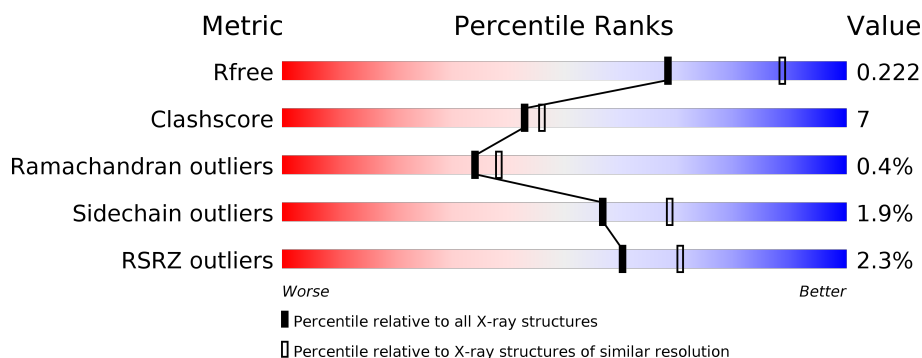
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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}	130704	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

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 on 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	814	<div> <div>2%</div> <div> <div></div> <div>82%</div> <div>14%</div> <div>.</div> </div> </div>
1	B	814	<div> <div>3%</div> <div> <div></div> <div>80%</div> <div>15%</div> <div>..</div> </div> </div>

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 13473 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 hydrolases family 2, sugar binding domain protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	781	Total	C	N	O	S	0	1	0
			6318	3977	1131	1181	29			
1	B	779	Total	C	N	O	S	0	1	0
			6295	3963	1128	1175	29			

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

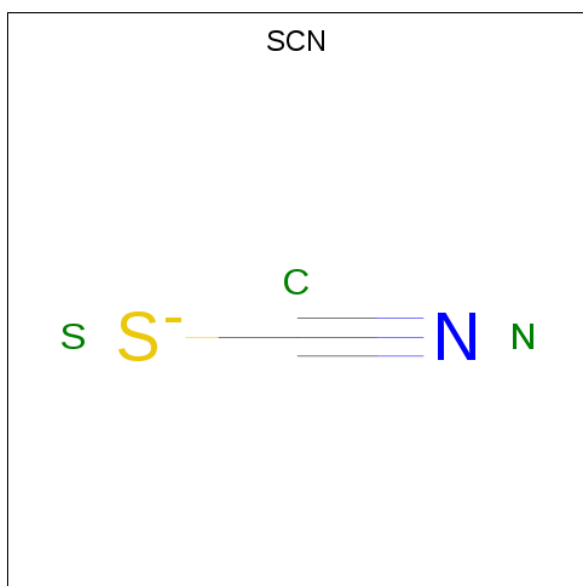
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	3	Total	Na	0	0
			3	3		
2	A	2	Total	Na	0	0
			2	2		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

- Molecule 4 is THIOCYANATE ION (three-letter code: SCN) (formula: CNS).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C N S 3 1 1 1	0	0
4	A	1	Total C N S 3 1 1 1	0	0

- Molecule 5 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	2	Total K 2 2	0	0
5	A	2	Total K 2 2	0	0

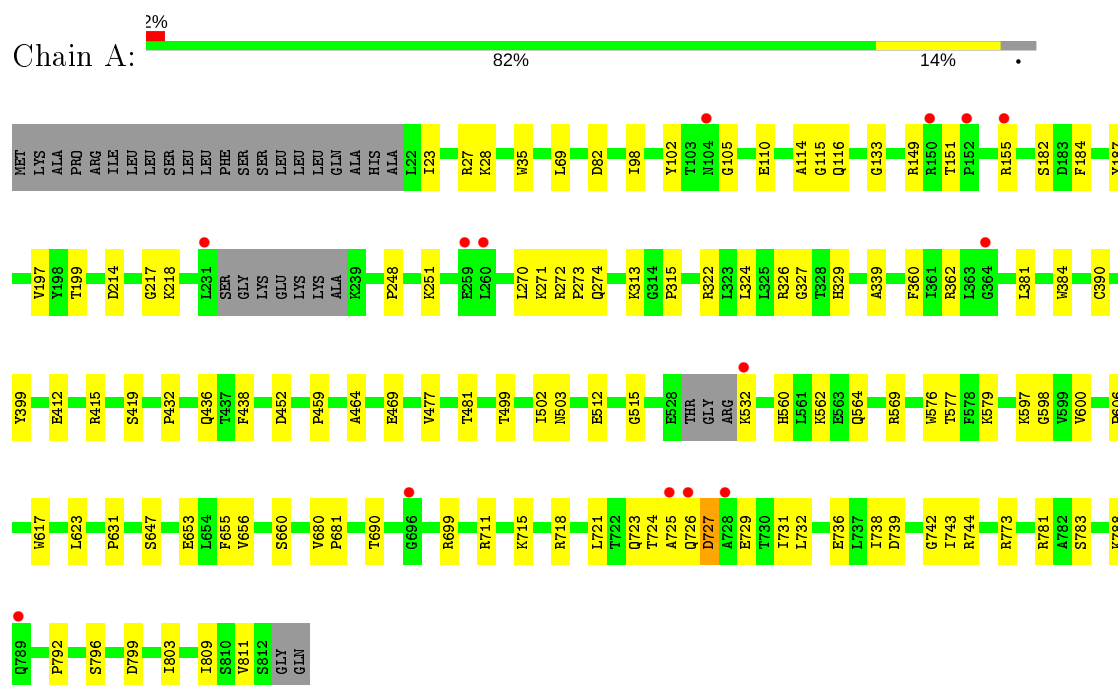
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	430	Total 430	O 430	0	0
6	B	385	Total 385	O 385	0	0

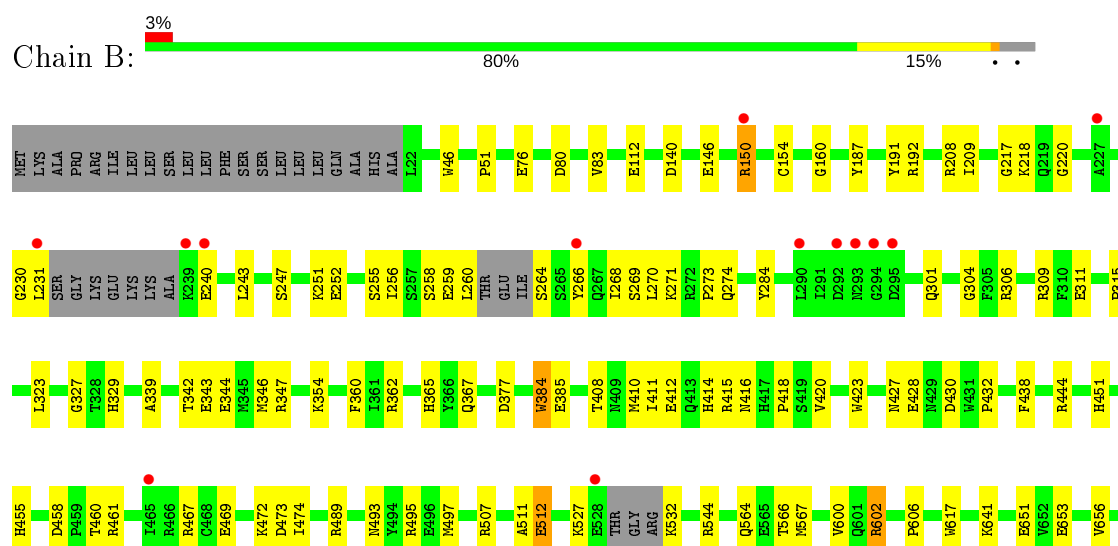
3 Residue-property plots [i](#)

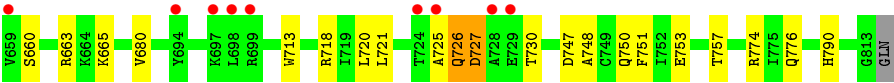
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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 hydrolases family 2, sugar binding domain protein



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4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	209.72Å 68.21Å 167.30Å 90.00° 122.49° 90.00°	Depositor
Resolution (Å)	29.09 – 2.35 29.09 – 2.35	Depositor EDS
% Data completeness (in resolution range)	100.0 (29.09-2.35) 100.0 (29.09-2.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.77 (at 2.36Å)	Xtriage
Refinement program	PHENIX (1.14_3260)	Depositor
R, R_{free}	0.175 , 0.222 0.175 , 0.222	Depositor DCC
R_{free} test set	2000 reflections (2.40%)	wwPDB-VP
Wilson B-factor (Å ²)	31.9	Xtriage
Anisotropy	0.966	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13473	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.54% 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, K, SCN, NA

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	0.44	0/6476	0.59	0/8767
1	B	0.46	0/6455	0.61	2/8736 (0.0%)
All	All	0.45	0/12931	0.60	2/17503 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	602	ARG	NE-CZ-NH2	-7.41	116.59	120.30
1	B	602	ARG	NE-CZ-NH1	6.24	123.42	120.30

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	6318	0	6097	68	0
1	B	6295	0	6075	102	0
2	A	2	0	0	0	0
2	B	3	0	0	0	0
3	A	18	0	24	1	0
3	B	12	0	16	3	0
4	A	6	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	2	0	0	0	0
5	B	2	0	0	0	0
6	A	430	0	0	18	1
6	B	385	0	0	24	2
All	All	13473	0	12212	169	2

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 (169) 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:154:CYS:SG	6:B:1171:HOH:O	2.23	0.95
1:A:459:PRO:O	6:A:1001:HOH:O	1.86	0.92
1:B:271:LYS:NZ	6:B:1005:HOH:O	2.11	0.83
1:A:452:ASP:OD2	6:A:1002:HOH:O	1.96	0.83
1:B:274:GLN:NE2	6:B:1007:HOH:O	2.13	0.81
1:B:231:LEU:O	6:B:1001:HOH:O	1.99	0.80
1:A:656:VAL:HG21	1:A:680:VAL:HG11	1.66	0.78
1:B:112:GLU:OE2	6:B:1002:HOH:O	2.02	0.76
1:A:792:PRO:O	6:A:1003:HOH:O	2.02	0.76
1:B:140:ASP:O	6:B:1003:HOH:O	2.04	0.76
1:B:726:GLN:HG3	1:B:727:ASP:HB2	1.69	0.74
1:B:146:GLU:OE2	1:B:150:ARG:NH2	2.20	0.74
1:B:264:SER:N	6:B:1014:HOH:O	2.21	0.73
1:B:495:ARG:NH2	1:B:566:THR:OG1	2.24	0.71
1:B:230:GLY:O	1:B:231:LEU:HB2	1.89	0.70
1:B:656:VAL:HG21	1:B:680:VAL:HG21	1.74	0.70
1:B:790:HIS:O	6:B:1004:HOH:O	2.09	0.69
3:A:904:GOL:H12	3:A:905:GOL:H12	1.74	0.68
1:B:208:ARG:HD2	6:B:1037:HOH:O	1.93	0.68
1:A:412:GLU:OE1	6:A:1005:HOH:O	2.12	0.68
1:A:214:ASP:OD2	1:A:218:LYS:HE2	1.95	0.67
1:A:23:ILE:HG21	1:A:199:THR:HG21	1.77	0.67
1:B:76:GLU:OE1	6:B:1006:HOH:O	2.13	0.67
1:B:252:GLU:OE2	1:B:255:SER:OG	2.13	0.66
1:A:151:THR:O	6:A:1006:HOH:O	2.15	0.65
1:B:256:ILE:HG12	1:B:266:TYR:OH	1.98	0.64
1:A:82:ASP:OD2	6:A:1007:HOH:O	2.15	0.64
1:A:503:ASN:OD1	1:A:569:ARG:NH2	2.32	0.63
1:B:713:TRP:HE1	1:B:747:ASP:HB2	1.62	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:323:LEU:O	6:B:1008:HOH:O	2.15	0.63
1:A:727:ASP:O	1:A:788:LYS:NZ	2.18	0.63
1:A:502:ILE:O	6:A:1008:HOH:O	2.16	0.63
1:B:51:PRO:O	6:B:1010:HOH:O	2.16	0.62
1:B:512:GLU:OE2	3:B:904:GOL:H31	2.00	0.62
1:A:725:ALA:HB2	1:A:731:ILE:HG23	1.84	0.60
1:B:469:GLU:OE1	6:B:1011:HOH:O	2.17	0.60
1:B:473:ASP:O	6:B:1009:HOH:O	2.16	0.60
1:A:736:GLU:OE2	1:A:744:ARG:NH2	2.35	0.59
1:A:98:ILE:O	1:A:149:ARG:NH2	2.32	0.59
1:A:781:ARG:NH2	6:A:1017:HOH:O	2.36	0.58
1:A:738:ILE:HD12	1:A:742:GLY:HA2	1.86	0.57
1:B:266:TYR:HE1	1:B:268:ILE:HG13	1.69	0.57
1:B:430:ASP:OD1	1:B:467:ARG:NH1	2.38	0.56
1:B:304:GLY:N	1:B:418:PRO:HG3	2.22	0.54
1:B:344:GLU:CD	1:B:344:GLU:H	2.12	0.54
1:B:347:ARG:HB3	1:B:347:ARG:HH21	1.73	0.54
1:B:420:VAL:HG11	1:B:423:TRP:CH2	2.43	0.53
1:B:415:ARG:NH1	6:B:1020:HOH:O	2.25	0.53
1:B:444:ARG:HG3	1:B:474:ILE:HD11	1.90	0.53
1:B:220:GLY:O	1:B:270:LEU:N	2.35	0.53
1:B:726:GLN:HG2	1:B:730:THR:HG22	1.90	0.53
1:A:329:HIS:CD2	1:A:362:ARG:HB3	2.44	0.53
1:B:266:TYR:CE1	1:B:268:ILE:HG13	2.43	0.53
1:B:385:GLU:OE1	1:B:414:HIS:ND1	2.31	0.53
1:B:460:THR:HG22	1:B:461:ARG:HG3	1.89	0.53
1:B:751:PHE:CE1	1:B:776:GLN:HG3	2.44	0.53
1:A:718:ARG:NH2	6:A:1004:HOH:O	2.11	0.53
1:B:256:ILE:HG12	1:B:266:TYR:CZ	2.43	0.53
1:A:432:PRO:HB3	1:A:438:PHE:CD1	2.44	0.52
1:B:653:GLU:OE2	1:B:660:SER:OG	2.23	0.52
1:B:651:GLU:HG2	1:B:665:LYS:HG2	1.90	0.52
1:B:343:GLU:O	1:B:347:ARG:HG3	2.10	0.51
1:B:713:TRP:NE1	1:B:747:ASP:HB2	2.24	0.51
1:B:458:ASP:OD1	1:B:460:THR:HB	2.11	0.51
1:A:796:SER:HB3	1:B:757:THR:HB	1.93	0.51
1:A:248:PRO:HG2	1:A:274:GLN:OE1	2.11	0.51
1:A:725:ALA:CB	1:A:731:ILE:HA	2.41	0.50
1:B:269:SER:O	1:B:270:LEU:HD23	2.11	0.50
1:B:354:LYS:NZ	6:B:1019:HOH:O	2.25	0.50
1:B:415:ARG:HD3	6:B:1020:HOH:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:329:HIS:CD2	1:B:362:ARG:HB3	2.46	0.50
1:A:28:LYS:HB3	1:A:197:VAL:HG22	1.92	0.50
1:B:258:SER:O	1:B:259:GLU:HG2	2.11	0.50
1:A:381:LEU:O	1:A:419:SER:HB2	2.12	0.49
1:B:377:ASP:CG	1:B:418:PRO:HD2	2.33	0.49
1:A:415:ARG:NH1	6:A:1015:HOH:O	2.33	0.49
1:A:322:ARG:NH2	1:A:324:LEU:HD12	2.28	0.49
1:A:732:LEU:HD11	1:A:783:SER:HB3	1.95	0.49
1:B:266:TYR:OH	1:B:268:ILE:HD11	2.13	0.49
1:B:385:GLU:HG2	1:B:410:MET:SD	2.53	0.49
1:B:564:GLN:HA	1:B:567:MET:CE	2.43	0.49
1:B:600:VAL:HG22	1:B:606:PRO:HA	1.94	0.49
1:A:792:PRO:HA	1:A:811:VAL:O	2.13	0.48
1:B:415:ARG:HG3	6:B:1015:HOH:O	2.13	0.48
1:B:365:HIS:CE1	3:B:904:GOL:H32	2.49	0.48
1:A:327:GLY:HA2	1:A:360:PHE:O	2.14	0.48
1:A:313:LYS:NZ	1:A:503:ASN:O	2.47	0.47
1:B:493:ASN:O	1:B:497:MET:HE2	2.13	0.47
1:B:564:GLN:HA	1:B:567:MET:HE2	1.96	0.47
1:A:711:ARG:NH2	6:A:1014:HOH:O	2.31	0.47
1:B:602:ARG:NH2	6:B:1043:HOH:O	2.46	0.47
1:B:247:SER:HA	1:B:284:TYR:CD1	2.49	0.47
1:A:436:GLN:HG2	1:B:46:TRP:CE3	2.50	0.47
1:A:680:VAL:HG12	1:A:681:PRO:O	2.15	0.47
1:A:655:PHE:HB2	1:A:690:THR:OG1	2.15	0.47
1:B:651:GLU:OE2	1:B:663:ARG:NE	2.41	0.47
1:B:726:GLN:CG	1:B:727:ASP:HB2	2.42	0.47
1:B:160:GLY:HA2	6:B:1171:HOH:O	2.15	0.47
1:B:187:TYR:CD2	1:B:339:ALA:HB2	2.50	0.47
1:B:718:ARG:NH2	1:B:720:LEU:HD21	2.30	0.47
1:A:579:LYS:HG3	1:A:600:VAL:HG23	1.95	0.46
1:A:499:THR:HG23	1:A:569:ARG:HH12	1.80	0.46
1:B:342:THR:O	1:B:346:MET:HG3	2.16	0.46
1:B:362:ARG:NH2	1:B:427:ASN:OD1	2.49	0.46
1:B:365:HIS:NE2	3:B:904:GOL:H32	2.31	0.46
1:B:112:GLU:O	1:B:192:ARG:HD2	2.16	0.46
1:B:411:ILE:O	1:B:415:ARG:HB3	2.16	0.45
1:A:623:LEU:HD12	1:A:647:SER:HB2	1.98	0.45
1:A:326:ARG:HB3	1:A:617:TRP:CE2	2.51	0.45
1:B:420:VAL:HG11	1:B:423:TRP:CZ3	2.52	0.45
1:A:773:ARG:NE	6:A:1029:HOH:O	2.45	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:600:VAL:HG12	1:A:606:PRO:HA	1.99	0.45
1:B:430:ASP:O	1:B:432:PRO:HD3	2.17	0.45
1:A:102:TYR:HB2	1:A:105:GLY:HA3	1.99	0.45
1:A:115:GLY:HA2	1:A:116:GLN:HA	1.67	0.44
1:A:464:ALA:HA	1:A:477:VAL:O	2.17	0.44
1:B:301:GLN:OE1	1:B:416:ASN:ND2	2.49	0.44
1:A:725:ALA:HB2	1:A:731:ILE:HA	1.98	0.44
1:B:469:GLU:O	1:B:472:LYS:HD2	2.17	0.44
1:A:481:THR:HA	6:A:1184:HOH:O	2.17	0.44
1:B:208:ARG:HG3	1:B:415:ARG:HD3	2.00	0.44
1:B:218:LYS:HD3	1:B:218:LYS:HA	1.72	0.44
1:B:367:GLN:HG3	1:B:414:HIS:CE1	2.52	0.44
1:A:23:ILE:N	1:A:110:GLU:OE2	2.44	0.44
1:B:427:ASN:ND2	1:B:428:GLU:HG3	2.33	0.44
1:A:739:ASP:OD2	1:A:743:ILE:HB	2.18	0.43
1:A:781:ARG:HD3	6:A:1115:HOH:O	2.18	0.43
1:B:507:ARG:NH2	6:B:1031:HOH:O	2.33	0.43
1:A:576:TRP:HA	1:A:577:THR:HA	1.84	0.43
1:B:309:ARG:NH1	1:B:311:GLU:HG3	2.34	0.43
1:A:469:GLU:HB3	6:A:1079:HOH:O	2.17	0.43
1:B:243:LEU:O	1:B:256:ILE:HD12	2.19	0.43
1:B:347:ARG:NH2	1:B:347:ARG:HB3	2.34	0.43
1:B:451:HIS:CE1	1:B:455:HIS:CD2	3.07	0.43
1:A:799:ASP:OD2	1:A:803:ILE:HG12	2.18	0.42
1:A:721:LEU:HD13	1:A:809:ILE:HG22	2.01	0.42
1:B:217:GLY:O	1:B:273:PRO:HD2	2.19	0.42
1:A:114:ALA:O	1:A:133:GLY:HA2	2.19	0.42
1:A:515:GLY:C	1:A:597:LYS:HG2	2.40	0.42
1:A:560:HIS:O	1:A:564:GLN:HG3	2.19	0.42
1:B:327:GLY:HA2	1:B:360:PHE:O	2.20	0.42
1:B:617:TRP:O	6:B:1013:HOH:O	2.20	0.42
1:A:270:LEU:CD2	1:A:273:PRO:HB3	2.50	0.42
1:A:217:GLY:O	1:A:272:ARG:HA	2.20	0.42
1:B:753:GLU:OE2	1:B:774:ARG:NE	2.47	0.42
1:B:209:ILE:O	6:B:1015:HOH:O	2.22	0.42
1:B:384:TRP:CD1	1:B:384:TRP:C	2.93	0.42
1:B:721:LEU:HA	1:B:721:LEU:HD23	1.86	0.42
1:B:80:ASP:HB3	1:B:83:VAL:HG22	2.02	0.42
1:B:432:PRO:HB3	1:B:438:PHE:CD1	2.55	0.41
1:A:577:THR:O	1:A:598:GLY:HA2	2.19	0.41
1:B:260:LEU:HD22	1:B:260:LEU:H	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:270:LEU:HD23	1:A:273:PRO:HB3	2.02	0.41
1:A:35:TRP:CD1	1:A:69:LEU:HG	2.56	0.41
1:A:631:PRO:HD2	6:A:1225:HOH:O	2.21	0.41
1:B:511:ALA:O	1:B:512:GLU:HB2	2.21	0.41
1:B:247:SER:HA	1:B:284:TYR:CE1	2.55	0.41
1:B:489:ARG:HD2	6:B:1100:HOH:O	2.20	0.41
1:B:493:ASN:C	1:B:497:MET:HE2	2.41	0.41
1:B:641:LYS:HA	1:B:641:LYS:HD2	1.82	0.41
1:A:182:SER:HB3	1:A:184:PHE:CE1	2.56	0.41
1:A:653:GLU:OE1	1:A:660:SER:OG	2.23	0.40
1:A:390:CYS:HB2	6:A:1272:HOH:O	2.21	0.40
1:A:773:ARG:HD3	6:A:1183:HOH:O	2.21	0.40
1:B:220:GLY:N	1:B:270:LEU:O	2.31	0.40
1:B:408:THR:O	1:B:412:GLU:HG3	2.22	0.40
1:A:736:GLU:OE2	1:A:744:ARG:NE	2.48	0.40
1:A:187:TYR:CD2	1:A:339:ALA:HB2	2.56	0.40
1:B:748:ALA:HB1	1:B:750:GLN:OE1	2.22	0.40

All (2) 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)
6:A:1338:HOH:O	6:B:1042:HOH:O[3_555]	1.66	0.54
6:B:1014:HOH:O	6:B:1359:HOH:O[2_555]	2.05	0.15

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	776/814 (95%)	743 (96%)	31 (4%)	2 (0%)	41 47
1	B	772/814 (95%)	733 (95%)	35 (4%)	4 (0%)	29 32

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1548/1628 (95%)	1476 (95%)	66 (4%)	6 (0%)	34 38

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	725	ALA
1	A	512	GLU
1	B	315	PRO
1	B	527	LYS
1	B	512	GLU
1	A	315	PRO

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	673/699 (96%)	658 (98%)	15 (2%)	52 63
1	B	670/699 (96%)	660 (98%)	10 (2%)	65 76
All	All	1343/1398 (96%)	1318 (98%)	25 (2%)	57 68

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	ARG
1	A	155	ARG
1	A	251	LYS
1	A	271	LYS
1	A	384	TRP
1	A	399	TYR
1	A	532	LYS
1	A	562	LYS
1	A	699	ARG
1	A	715	LYS
1	A	723	GLN
1	A	724	THR

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Mol	Chain	Res	Type
1	A	726	GLN
1	A	727	ASP
1	A	729	GLU
1	B	150	ARG
1	B	191	TYR
1	B	240	GLU
1	B	251	LYS
1	B	306	ARG
1	B	384	TRP
1	B	532	LYS
1	B	544	ARG
1	B	726	GLN
1	B	727	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	274	GLN
1	B	416	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 9 are monoatomic - leaving 7 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
4	SCN	A	906	-	1,2,2	1.01	0	0,1,1	0.00	-
3	GOL	A	904	-	5,5,5	0.70	0	5,5,5	0.94	0
3	GOL	B	904	-	5,5,5	0.75	0	5,5,5	1.19	1 (20%)
4	SCN	A	907	-	1,2,2	1.02	0	0,1,1	0.00	-
3	GOL	A	905	-	5,5,5	1.30	1 (20%)	5,5,5	0.85	0
3	GOL	A	903	-	5,5,5	0.80	0	5,5,5	1.07	0
3	GOL	B	905	-	5,5,5	1.25	1 (20%)	5,5,5	0.72	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	903	-	-	0/4/4/4	-
3	GOL	B	905	-	-	0/4/4/4	-
3	GOL	A	904	-	-	4/4/4/4	-
3	GOL	B	904	-	-	4/4/4/4	-
3	GOL	A	905	-	-	1/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	905	GOL	C3-C2	2.23	1.60	1.51
3	A	905	GOL	C1-C2	2.23	1.60	1.51

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	904	GOL	C3-C2-C1	-2.29	102.79	111.70

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	904	GOL	C1-C2-C3-O3
3	B	904	GOL	C1-C2-C3-O3
3	B	904	GOL	O2-C2-C3-O3
3	A	904	GOL	O2-C2-C3-O3
3	A	904	GOL	O1-C1-C2-C3
3	B	904	GOL	O1-C1-C2-C3
3	A	905	GOL	O2-C2-C3-O3
3	A	904	GOL	O1-C1-C2-O2
3	B	904	GOL	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	904	GOL	1	0
3	B	904	GOL	3	0
3	A	905	GOL	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 ⓘ

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	781/814 (95%)	-0.18	14 (1%) 68 77	18, 32, 58, 72	0
1	B	779/814 (95%)	-0.13	22 (2%) 53 64	19, 33, 54, 77	0
All	All	1560/1628 (95%)	-0.15	36 (2%) 60 70	18, 33, 56, 77	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	726	GLN	7.1
1	A	725	ALA	4.7
1	B	728	ALA	4.1
1	B	239	LYS	3.8
1	B	698	LEU	3.8
1	B	293	ASN	3.7
1	A	155	ARG	3.4
1	A	728	ALA	3.4
1	B	231	LEU	3.4
1	B	150	ARG	3.3
1	B	697	LYS	3.3
1	B	295	ASP	3.1
1	B	699	ARG	3.1
1	B	240	GLU	2.7
1	B	729	GLU	2.7
1	A	104	ASN	2.6
1	A	259	GLU	2.6
1	A	789	GLN	2.5
1	B	292	ASP	2.5
1	B	227	ALA	2.4
1	B	725	ALA	2.4
1	B	266	TYR	2.3
1	B	465	ILE	2.3
1	B	528	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	231	LEU	2.2
1	A	696	GLY	2.2
1	A	364	GLY	2.2
1	B	290	LEU	2.2
1	A	152	PRO	2.2
1	B	294	GLY	2.2
1	A	532	LYS	2.1
1	B	694	TYR	2.1
1	B	659	VAL	2.1
1	A	150	ARG	2.0
1	A	260	LEU	2.0
1	B	724	THR	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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
2	NA	B	902	1/1	0.86	0.27	36,36,36,36	0
4	SCN	A	907	3/3	0.88	0.16	35,35,36,37	0
4	SCN	A	906	3/3	0.89	0.24	47,47,51,59	0
3	GOL	A	905	6/6	0.91	0.19	20,30,31,32	0
2	NA	B	901	1/1	0.92	0.17	38,38,38,38	0
2	NA	B	903	1/1	0.93	0.50	47,47,47,47	0
3	GOL	B	904	6/6	0.93	0.23	31,33,36,38	0
3	GOL	B	905	6/6	0.96	0.10	37,38,39,45	0
3	GOL	A	903	6/6	0.97	0.10	36,37,41,41	0
3	GOL	A	904	6/6	0.98	0.12	20,24,25,27	0
2	NA	A	902	1/1	0.98	0.07	25,25,25,25	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	K	B	906	1/1	0.99	0.06	23,23,23,23	0
5	K	A	908	1/1	0.99	0.05	20,20,20,20	0
2	NA	A	901	1/1	0.99	0.06	30,30,30,30	0
5	K	A	909	1/1	0.99	0.09	21,21,21,21	0
5	K	B	907	1/1	0.99	0.06	27,27,27,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.