



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 10, 2017 – 04:59 AM EDT

PDB ID : 2FHF  
Title : Crystal Structure Analysis of Klebsiella pneumoniae pullulanase complexed with maltotetraose  
Authors : Mikami, B.; Iwamoto, H.; Katsuya, Y.; Yoon, H.-J.; Demirkan-Sarikaya, E.; Malle, D.  
Deposited on : unknown  
Resolution : 1.65 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20030345  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20030345

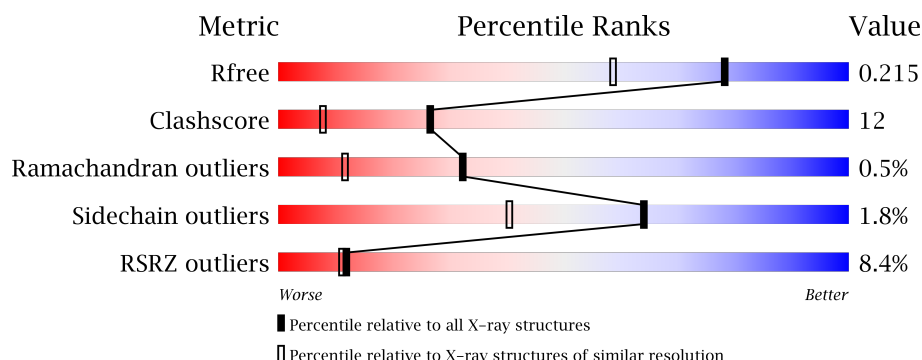
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.65 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))
$R_{free}$	100719	1368 (1.66-1.66)
Clashscore	112137	1468 (1.66-1.66)
Ramachandran outliers	110173	1438 (1.66-1.66)
Sidechain outliers	110143	1438 (1.66-1.66)
RSRZ outliers	101464	1371 (1.66-1.66)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	1083	<div> <div>8%</div> <div>79%</div> <div>17%</div> <div>..</div> </div>

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
2	GLC	A	1091	-	-	-	X
2	GLC	A	1092	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GLC	A	1099	-	-	-	X

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9441 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.

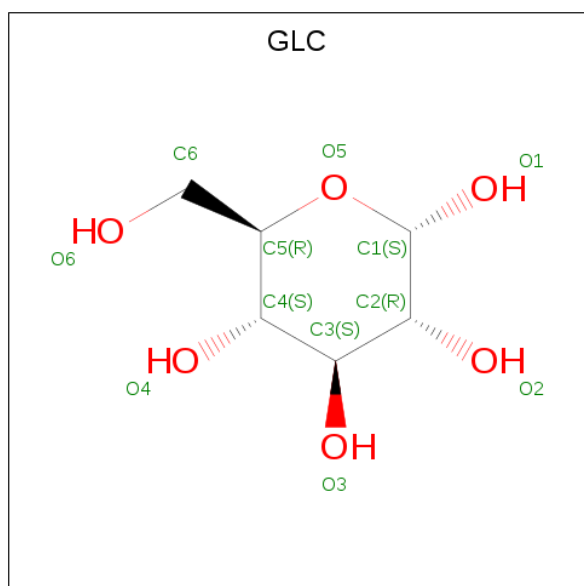
- Molecule 1 is a protein called pullulanase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	1052	8126	5081	1382	1636	27	0	18	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	680	LEU	GLY	ENGINEERED	GB 149300
A	882	LEU	VAL	ENGINEERED	GB 149300

- Molecule 2 is ALPHA-D-GLUCOSE (three-letter code: GLC) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			11	6	5		
2	A	1	Total	C	O	0	0
			11	6	5		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			11	6	5		
2	A	1	Total	C	O	0	0
			12	6	6		
2	A	1	Total	C	O	0	0
			11	6	5		
2	A	1	Total	C	O	0	0
			11	6	5		
2	A	1	Total	C	O	0	0
			11	6	5		
2	A	1	Total	C	O	0	0
			12	6	6		
2	A	1	Total	C	O	0	0
			11	6	5		
2	A	1	Total	C	O	0	0
			12	6	6		

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	5	Total	Ca	0	0
			5	5		

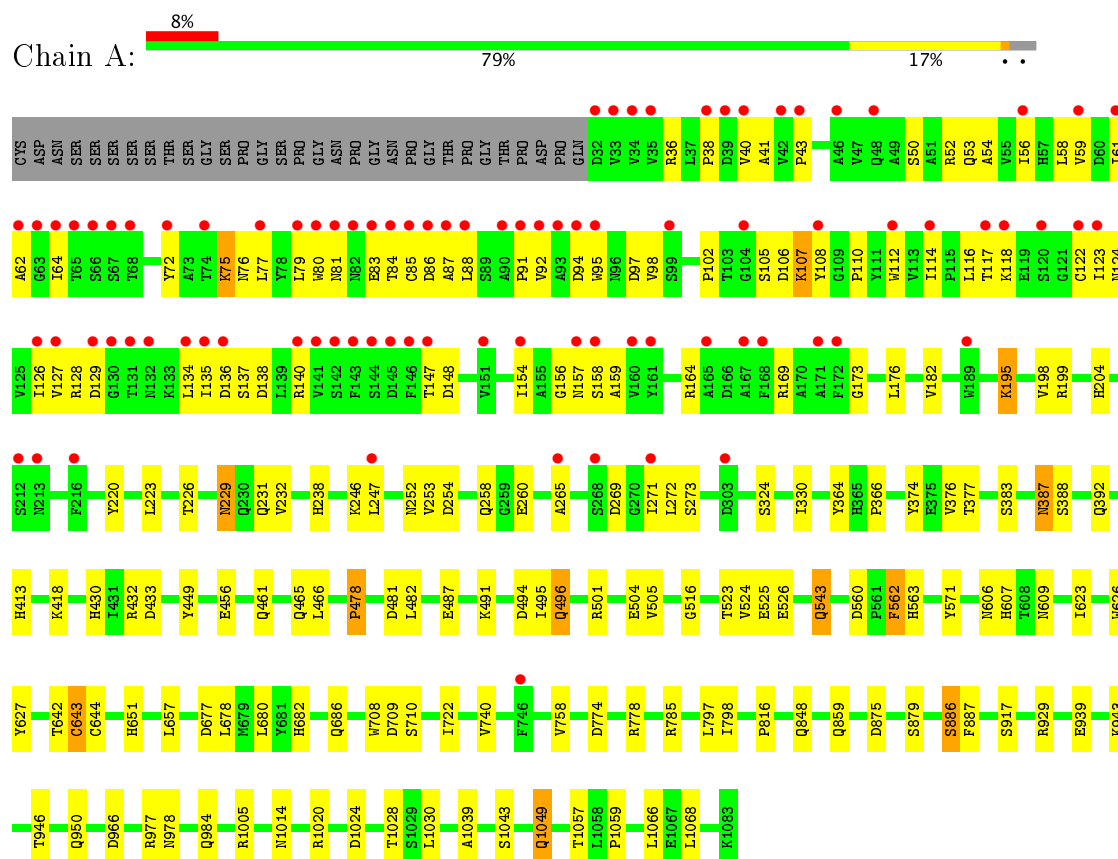
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1197	Total	O	0	0
			1197	1197		

### 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: pullulanase



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	150.16Å 60.51Å 134.66Å 90.00° 111.87° 90.00°	Depositor
Resolution (Å)	14.98 – 1.65 73.63 – 1.60	Depositor EDS
% Data completeness (in resolution range)	85.7 (14.98-1.65) 84.5 (73.63-1.60)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.66 (at 1.60Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.178 , 0.206 0.188 , 0.215	Depositor DCC
$R_{free}$ test set	11574 reflections (10.01%)	DCC
Wilson B-factor (Å <sup>2</sup> )	18.2	Xtriage
Anisotropy	0.500	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 54.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	9441	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.06% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

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

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

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

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	0.30	0/8363	0.60	0/11388

There are no bond length outliers.

There are no bond angle outliers.

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	8126	0	7830	186	0
2	A	113	0	99	6	0
3	A	5	0	0	0	0
4	A	1197	0	0	29	0
All	All	9441	0	7929	188	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 12.

All (188) 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:740:VAL:HA	1:A:758:VAL:HG12	1.44	0.97
1:A:81:ASN:HB2	1:A:86:ASP:HA	1.54	0.89
1:A:606:ASN:HD21	1:A:607:HIS:HD2	1.20	0.88
1:A:978:ASN:HD21	1:A:984:GLN:H	1.20	0.88
2:A:1093:GLC:H62	2:A:1098:GLC:O1	1.77	0.85
1:A:722:ILE:HD11	4:A:1973:HOH:O	1.78	0.83
1:A:677:ASP:OD1	2:A:1098:GLC:H1	1.80	0.82
1:A:81:ASN:HB3	1:A:88:LEU:HD13	1.62	0.80
1:A:680:LEU:HD11	4:A:1539:HOH:O	1.82	0.80
1:A:195:LYS:HG2	1:A:265:ALA:HB1	1.63	0.79
1:A:977:ARG:NH1	1:A:1024:ASP:HB3	1.98	0.79
1:A:722:ILE:HG12	4:A:1660:HOH:O	1.84	0.78
1:A:79:LEU:HD11	1:A:114:ILE:HD12	1.66	0.77
1:A:606:ASN:ND2	1:A:607:HIS:HD2	1.82	0.77
1:A:516:GLY:HA3	4:A:1195:HOH:O	1.85	0.77
1:A:324:SER:HB3	1:A:330[B]:ILE:HD11	1.68	0.76
1:A:157:ASN:HD21	1:A:159:ALA:HB3	1.50	0.75
1:A:229:ASN:ND2	1:A:232:VAL:H	1.88	0.72
1:A:501:ARG:O	1:A:504:GLU:HG2	1.89	0.72
1:A:560:ASP:HB3	1:A:609:ASN:ND2	2.03	0.72
1:A:680:LEU:HG	1:A:710:SER:HB3	1.73	0.71
1:A:560:ASP:HB3	1:A:609:ASN:HD22	1.57	0.70
1:A:376:VAL:HB	1:A:623[B]:ILE:HD11	1.74	0.69
1:A:88:LEU:HD12	1:A:88:LEU:H	1.58	0.68
1:A:229:ASN:HD21	1:A:232:VAL:HG23	1.57	0.68
1:A:740:VAL:HA	1:A:758:VAL:CG1	2.20	0.67
1:A:1030:LEU:HG	1:A:1066:LEU:HB3	1.76	0.67
1:A:505[A]:VAL:HG22	4:A:1366:HOH:O	1.95	0.66
1:A:107:LYS:HE3	1:A:107:LYS:H	1.61	0.66
1:A:627:TYR:O	1:A:651:HIS:HD2	1.79	0.66
1:A:157:ASN:ND2	1:A:159:ALA:HB3	2.11	0.65
1:A:84:THR:HG21	1:A:140:ARG:HH11	1.61	0.65
1:A:977:ARG:HH11	1:A:1024:ASP:HB3	1.60	0.65
1:A:680:LEU:HD13	1:A:708:TRP:HB2	1.78	0.64
1:A:76:ASN:CG	1:A:98:VAL:HG23	2.19	0.63
1:A:59:VAL:HB	1:A:154:ILE:HG12	1.81	0.62
1:A:1005:ARG:HB3	1:A:1005:ARG:NH1	2.14	0.62
1:A:677:ASP:OD1	2:A:1098:GLC:C1	2.47	0.62
1:A:127:VAL:HB	1:A:135:ILE:HD13	1.81	0.62
1:A:127:VAL:HG12	1:A:134:LEU:HD12	1.81	0.61
1:A:83:GLU:HG3	1:A:84:THR:N	2.16	0.60
1:A:229:ASN:HD21	1:A:232:VAL:CG2	2.15	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:246:LYS:HD3	1:A:247:LEU:O	2.02	0.59
1:A:682:HIS:HD2	1:A:686:GLN:NE2	2.01	0.58
1:A:1028:THR:HG22	1:A:1068:LEU:HD21	1.86	0.58
1:A:495:ILE:HA	1:A:524:VAL:HB	1.86	0.58
1:A:129:ASP:HB2	1:A:134:LEU:HD21	1.86	0.57
1:A:816:PRO:HG2	4:A:1973:HOH:O	2.03	0.57
1:A:84:THR:HG21	1:A:140:ARG:NH1	2.19	0.57
1:A:680:LEU:N	1:A:680:LEU:HD23	2.19	0.57
1:A:606:ASN:ND2	1:A:607:HIS:CD2	2.68	0.57
1:A:157:ASN:HB3	4:A:2296:HOH:O	2.05	0.57
1:A:722:ILE:HD12	4:A:1974:HOH:O	2.05	0.57
1:A:98:VAL:O	1:A:98:VAL:HG22	2.05	0.57
1:A:56:ILE:CG2	1:A:112:TRP:HB2	2.35	0.56
1:A:83:GLU:HG3	1:A:84:THR:H	1.69	0.56
1:A:680:LEU:HD23	1:A:680:LEU:H	1.69	0.56
1:A:1049:GLN:HG3	1:A:1057:THR:HB	1.87	0.56
1:A:682:HIS:HD2	1:A:686:GLN:HE22	1.53	0.56
1:A:126:ILE:HD11	1:A:138:ASP:OD1	2.05	0.56
1:A:229:ASN:HD21	1:A:232:VAL:H	1.51	0.56
1:A:642:THR:O	1:A:643:CYS:HB3	2.06	0.56
1:A:466:LEU:HD23	1:A:950:GLN:HE21	1.71	0.55
1:A:108:TYR:HB3	4:A:2279:HOH:O	2.05	0.55
1:A:56:ILE:HG22	1:A:112:TRP:HB2	1.87	0.55
1:A:198:VAL:HG12	1:A:223:LEU:HD12	1.89	0.55
1:A:774:ASP:HB3	4:A:2292:HOH:O	2.06	0.54
1:A:87:ALA:O	1:A:117:THR:HG22	2.07	0.54
1:A:491:LYS:C	1:A:505[A]:VAL:HG21	2.28	0.54
1:A:254:ASP:HB2	4:A:2122:HOH:O	2.07	0.54
1:A:387:ASN:ND2	1:A:487:GLU:H	2.05	0.54
1:A:106:ASP:HB2	1:A:107:LYS:NZ	2.23	0.54
1:A:1005:ARG:HB3	1:A:1005:ARG:HH11	1.73	0.54
1:A:1039:ALA:HB3	1:A:1043:SER:HB2	1.90	0.53
1:A:496:GLN:NE2	1:A:496:GLN:H	2.05	0.53
1:A:134:LEU:C	1:A:135:ILE:HD12	2.28	0.53
1:A:430:HIS:CD2	1:A:432:ARG:H	2.27	0.53
1:A:946:THR:O	1:A:950:GLN:HG3	2.09	0.53
1:A:465:GLN:HG3	1:A:950:GLN:HE22	1.75	0.52
1:A:680:LEU:HD12	1:A:710:SER:N	2.24	0.52
1:A:62:ALA:HB3	4:A:2256:HOH:O	2.10	0.51
1:A:79:LEU:HB3	1:A:88:LEU:HD23	1.91	0.51
1:A:43:PRO:HB3	1:A:108:TYR:CG	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:682:HIS:CD2	1:A:686:GLN:HE22	2.29	0.51
1:A:40:VAL:HG13	1:A:41:ALA:N	2.26	0.51
1:A:229:ASN:HD21	1:A:232:VAL:CB	2.24	0.50
1:A:680:LEU:CD2	1:A:680:LEU:H	2.25	0.50
1:A:107:LYS:CE	1:A:107:LYS:H	2.24	0.50
1:A:642:THR:O	1:A:643:CYS:CB	2.60	0.50
1:A:85:CYS:HB3	1:A:122:CYS:C	2.32	0.49
1:A:54:ALA:HB2	1:A:116:LEU:HD11	1.94	0.49
1:A:38:PRO:O	1:A:169:ARG:HD3	2.12	0.49
1:A:85:CYS:HB3	1:A:122:CYS:O	2.12	0.49
1:A:61:ILE:O	1:A:61:ILE:HG13	2.12	0.49
1:A:117:THR:HG23	1:A:118:LYS:HG2	1.94	0.49
1:A:80:TRP:O	1:A:124:ASN:HB2	2.12	0.49
1:A:523:THR:OG1	1:A:526:GLU:HG3	2.13	0.48
1:A:204:HIS:HE1	4:A:1154:HOH:O	1.97	0.48
1:A:76:ASN:ND2	1:A:128:ARG:HH21	2.11	0.48
1:A:182:VAL:HG21	1:A:232:VAL:HG11	1.96	0.48
1:A:81:ASN:ND2	1:A:91:PRO:HG3	2.29	0.48
1:A:978:ASN:ND2	1:A:984:GLN:H	2.01	0.48
1:A:680:LEU:HD12	1:A:709:ASP:C	2.35	0.48
1:A:88:LEU:HD12	1:A:88:LEU:N	2.26	0.48
1:A:1028:THR:HG22	1:A:1068:LEU:CD2	2.43	0.47
1:A:107:LYS:HD3	1:A:108:TYR:CE2	2.48	0.47
1:A:1066:LEU:N	1:A:1066:LEU:HD22	2.28	0.47
1:A:88:LEU:HD11	1:A:123:ILE:HA	1.95	0.47
1:A:740:VAL:HG22	1:A:758:VAL:CG1	2.45	0.47
1:A:678:LEU:C	1:A:680:LEU:HD23	2.35	0.47
1:A:176:LEU:HA	4:A:1715:HOH:O	2.14	0.47
1:A:72:TYR:CZ	1:A:110:PRO:HD3	2.50	0.47
1:A:157:ASN:O	1:A:158:SER:HB3	2.15	0.47
1:A:430:HIS:HD2	1:A:433:ASP:H	1.62	0.47
1:A:461[A]:GLN:NE2	4:A:1931:HOH:O	2.48	0.47
1:A:1014:ASN:HD21	1:A:1020:ARG:HH11	1.61	0.46
1:A:778:ARG:HD2	4:A:1519:HOH:O	2.14	0.46
1:A:253:VAL:HG23	1:A:254:ASP:N	2.29	0.46
1:A:105:SER:HB3	4:A:2282:HOH:O	2.14	0.46
1:A:260:GLU:HB2	1:A:364:TYR:CE1	2.51	0.46
1:A:107:LYS:HD3	1:A:108:TYR:CD2	2.51	0.46
1:A:238:HIS:CD2	1:A:238:HIS:H	2.33	0.46
1:A:92:VAL:HG21	1:A:97:ASP:OD2	2.16	0.46
1:A:147:THR:HG23	1:A:148:ASP:N	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:465:GLN:HG3	1:A:950:GLN:NE2	2.30	0.45
1:A:95:TRP:CZ3	2:A:1100:GLC:H2	2.51	0.45
1:A:229:ASN:HD22	1:A:229:ASN:C	2.17	0.45
1:A:1049:GLN:CG	1:A:1057:THR:HB	2.47	0.45
1:A:272:LEU:C	1:A:272:LEU:HD23	2.36	0.45
1:A:59:VAL:HG13	4:A:2279:HOH:O	2.16	0.45
1:A:199:ARG:HD3	1:A:220:TYR:CE2	2.51	0.45
1:A:377:THR:OG1	1:A:563:HIS:HE1	1.99	0.45
1:A:525:GLU:HG2	4:A:2178:HOH:O	2.15	0.45
1:A:238:HIS:HE1	4:A:1433:HOH:O	1.99	0.45
1:A:383:SER:HB3	1:A:392:GLN:HB3	1.98	0.45
1:A:80:TRP:CZ2	1:A:124:ASN:HB3	2.52	0.44
1:A:366:PRO:HB2	1:A:626:TRP:CE2	2.53	0.44
1:A:491:LYS:O	1:A:505[A]:VAL:HG21	2.17	0.44
1:A:199:ARG:HD3	1:A:220:TYR:CD2	2.52	0.44
1:A:643:CYS:SG	1:A:644:CYS:N	2.91	0.44
1:A:798[B]:ILE:HD13	4:A:1401:HOH:O	2.18	0.44
1:A:449:TYR:CZ	1:A:571:TYR:HB2	2.53	0.43
1:A:785:ARG:NH2	4:A:1359:HOH:O	2.50	0.43
1:A:886:SER:O	1:A:887:PHE:HB2	2.19	0.43
1:A:40:VAL:HG13	1:A:41:ALA:H	1.83	0.43
1:A:740:VAL:HG22	1:A:758:VAL:HG11	2.00	0.43
1:A:229:ASN:ND2	1:A:232:VAL:HB	2.33	0.43
1:A:269:ASP:OD1	1:A:271:ILE:HG13	2.18	0.43
2:A:1093:GLC:H62	2:A:1098:GLC:HO1	1.81	0.43
1:A:135:ILE:N	1:A:135:ILE:HD12	2.34	0.43
1:A:272:LEU:HD23	1:A:273:SER:N	2.33	0.43
1:A:36:ARG:HG2	1:A:36:ARG:HH11	1.84	0.43
1:A:56:ILE:HG21	1:A:77:LEU:HD11	2.00	0.43
1:A:413:HIS:HD2	4:A:1172:HOH:O	2.02	0.42
1:A:418:LYS:HD2	1:A:966[B]:ASP:OD1	2.19	0.42
1:A:52:ARG:HH11	1:A:52:ARG:HG3	1.84	0.42
1:A:56:ILE:HG23	1:A:112:TRP:HE3	1.84	0.42
1:A:75:LYS:O	1:A:102:PRO:HD3	2.20	0.42
1:A:231:GLN:N	1:A:231:GLN:OE1	2.51	0.42
1:A:50:SER:H	1:A:53:GLN:NE2	2.18	0.42
1:A:465:GLN:CG	1:A:950:GLN:HE22	2.33	0.42
1:A:606:ASN:HD21	1:A:607:HIS:CD2	2.13	0.41
1:A:682:HIS:CD2	1:A:686:GLN:NE2	2.85	0.41
1:A:108:TYR:HB2	1:A:164:ARG:NH1	2.34	0.41
1:A:1005:ARG:HH11	1:A:1005:ARG:CB	2.34	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:657:LEU:C	1:A:657:LEU:HD23	2.40	0.41
1:A:95:TRP:CH2	2:A:1100:GLC:H2	2.56	0.41
1:A:106:ASP:HB2	1:A:107:LYS:HZ1	1.85	0.41
1:A:481:ASP:OD2	1:A:563:HIS:HD2	2.03	0.41
1:A:543:GLN:NE2	1:A:917:SER:H	2.19	0.41
1:A:939:GLU:OE1	1:A:943:LYS:HE3	2.21	0.41
1:A:797:LEU:HD12	1:A:797:LEU:C	2.40	0.41
1:A:110:PRO:HB3	4:A:2282:HOH:O	2.20	0.41
1:A:494:ASP:HB3	1:A:496:GLN:HE22	1.85	0.41
1:A:58:LEU:HB2	4:A:2283:HOH:O	2.21	0.41
1:A:875:ASP:OD1	1:A:879:SER:HB2	2.21	0.41
1:A:387:ASN:HD22	1:A:487:GLU:HG3	1.86	0.41
1:A:226:THR:HB	1:A:246:LYS:HB2	2.03	0.41
1:A:156:GLY:HA2	4:A:2222:HOH:O	2.20	0.41
1:A:929:ARG:HD2	4:A:1584:HOH:O	2.20	0.41
1:A:173:GLY:HA2	4:A:2189:HOH:O	2.21	0.41
1:A:848:GLN:HG3	4:A:1506:HOH:O	2.21	0.41
1:A:254:ASP:O	1:A:258:GLN:HG2	2.22	0.40
1:A:388:SER:HB2	1:A:562:PHE:CE1	2.55	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	1068/1083 (99%)	1036 (97%)	27 (2%)	5 (0%)	32 12

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	478	PRO
1	A	643	CYS

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Mol	Chain	Res	Type
1	A	136	ASP
1	A	94	ASP
1	A	137	SER

### 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	884/891 (99%)	868 (98%)	16 (2%)	64	40

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	64	ILE
1	A	75	LYS
1	A	107	LYS
1	A	195	LYS
1	A	229	ASN
1	A	252	ASN
1	A	374	TYR
1	A	387	ASN
1	A	478	PRO
1	A	482	LEU
1	A	496	GLN
1	A	543	GLN
1	A	562	PHE
1	A	859	GLN
1	A	886	SER
1	A	1049	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (51) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	53	GLN
1	A	57	HIS

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Mol	Chain	Res	Type
1	A	81	ASN
1	A	124	ASN
1	A	157	ASN
1	A	194	ASN
1	A	204	HIS
1	A	213	ASN
1	A	229	ASN
1	A	230	GLN
1	A	238	HIS
1	A	252	ASN
1	A	258	GLN
1	A	279	GLN
1	A	316	GLN
1	A	317	GLN
1	A	387	ASN
1	A	392	GLN
1	A	413	HIS
1	A	430	HIS
1	A	439	GLN
1	A	455	GLN
1	A	458	ASN
1	A	465	GLN
1	A	496	GLN
1	A	533	GLN
1	A	534	ASN
1	A	541	GLN
1	A	543	GLN
1	A	551	GLN
1	A	563	HIS
1	A	606	ASN
1	A	607	HIS
1	A	609	ASN
1	A	629	GLN
1	A	651	HIS
1	A	682	HIS
1	A	686	GLN
1	A	711	ASN
1	A	848	GLN
1	A	859	GLN
1	A	899	ASN
1	A	911	ASN
1	A	950	GLN

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Mol	Chain	Res	Type
1	A	978	ASN
1	A	983	GLN
1	A	1014	ASN
1	A	1023	GLN
1	A	1037	GLN
1	A	1049	GLN
1	A	1074	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 15 ligands modelled in this entry, 5 are monoatomic - leaving 10 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	$\# Z  > 2$
2	GLC	A	1091	2	11,11,12	0.46	0	13,15,17	0.63	0
2	GLC	A	1092	2	11,11,12	0.52	0	13,15,17	0.61	0
2	GLC	A	1093	2	11,11,12	0.51	0	13,15,17	0.71	0
2	GLC	A	1094	2	12,12,12	0.51	0	17,17,17	0.52	0
2	GLC	A	1095	2	11,11,12	0.43	0	13,15,17	0.55	0
2	GLC	A	1096	2	11,11,12	0.49	0	13,15,17	0.67	0
2	GLC	A	1097	2	11,11,12	0.52	0	13,15,17	0.73	1 (7%)



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	GLC	A	1098	2	12,12,12	0.42	0	17,17,17	0.60	0
2	GLC	A	1099	2	11,11,12	0.47	0	13,15,17	0.61	0
2	GLC	A	1100	2	12,12,12	0.41	0	17,17,17	0.31	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	GLC	A	1091	2	-	0/2/19/22	0/1/1/1
2	GLC	A	1092	2	-	0/2/19/22	0/1/1/1
2	GLC	A	1093	2	-	0/2/19/22	0/1/1/1
2	GLC	A	1094	2	-	0/2/22/22	0/1/1/1
2	GLC	A	1095	2	-	0/2/19/22	0/1/1/1
2	GLC	A	1096	2	-	0/2/19/22	0/1/1/1
2	GLC	A	1097	2	-	0/2/19/22	0/1/1/1
2	GLC	A	1098	2	-	0/2/22/22	0/1/1/1
2	GLC	A	1099	2	-	0/2/19/22	0/1/1/1
2	GLC	A	1100	2	-	0/2/22/22	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1097	GLC	C1-O5-C5	2.36	115.42	112.17

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1093	GLC	2	0
2	A	1098	GLC	4	0
2	A	1100	GLC	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	1052/1083 (97%)	0.05	88 (8%) <b>12</b> <b>11</b>	13, 21, 56, 74	124 (11%)

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	84	THR	9.5
1	A	83	GLU	8.1
1	A	81	ASN	7.9
1	A	93	ALA	7.5
1	A	120	SER	7.2
1	A	172	PHE	6.6
1	A	143	PHE	6.0
1	A	40	VAL	6.0
1	A	72	TYR	5.3
1	A	142	SER	5.3
1	A	129	ASP	5.3
1	A	132	ASN	5.2
1	A	117	THR	5.0
1	A	90	ALA	4.7
1	A	131	THR	4.7
1	A	147	THR	4.7
1	A	168	PHE	4.5
1	A	303	ASP	4.5
1	A	62	ALA	4.5
1	A	42	VAL	4.4
1	A	82	ASN	4.3
1	A	136	ASP	4.3
1	A	56	ILE	4.1
1	A	43	PRO	4.1
1	A	95	TRP	4.0
1	A	158	SER	4.0
1	A	122	CYS	4.0

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Mol	Chain	Res	Type	RSRZ
1	A	154	ILE	3.9
1	A	80	TRP	3.9
1	A	92	VAL	3.8
1	A	146	PHE	3.7
1	A	59	VAL	3.6
1	A	123	ILE	3.6
1	A	171	ALA	3.5
1	A	85	CYS	3.5
1	A	64	ILE	3.4
1	A	99	SER	3.4
1	A	88	LEU	3.3
1	A	79	LEU	3.2
1	A	135	ILE	3.2
1	A	118	LYS	3.2
1	A	271	ILE	3.1
1	A	35	VAL	3.1
1	A	145	ASP	3.0
1	A	86	ASP	3.0
1	A	108	TYR	3.0
1	A	74	THR	2.9
1	A	134	LEU	2.9
1	A	39	ASP	2.8
1	A	77	LEU	2.8
1	A	38	PRO	2.8
1	A	265	ALA	2.7
1	A	46	ALA	2.7
1	A	144	SER	2.7
1	A	140	ARG	2.7
1	A	161	TYR	2.7
1	A	141	VAL	2.6
1	A	65	THR	2.6
1	A	94	ASP	2.6
1	A	165	ALA	2.5
1	A	212	SER	2.5
1	A	67	SER	2.5
1	A	126	ILE	2.5
1	A	157	ASN	2.5
1	A	160	VAL	2.4
1	A	87	ALA	2.4
1	A	247	LEU	2.4
1	A	68	THR	2.4
1	A	63	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	32	ASP	2.4
1	A	213	ASN	2.4
1	A	66	SER	2.3
1	A	127	VAL	2.3
1	A	746	PHE	2.3
1	A	91	PRO	2.3
1	A	130	GLY	2.3
1	A	268	SER	2.2
1	A	114	ILE	2.2
1	A	216	PHE	2.2
1	A	34	VAL	2.2
1	A	104	GLY	2.2
1	A	112	TRP	2.2
1	A	61	ILE	2.1
1	A	189	TRP	2.1
1	A	48	GLN	2.1
1	A	167	ALA	2.1
1	A	151	VAL	2.1
1	A	33	VAL	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	GLC	A	1091	11/12	0.82	0.27	3.84	33,35,36,36	0
2	GLC	A	1092	11/12	0.81	0.25	2.57	27,31,33,34	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	GLC	A	1099	11/12	0.71	0.29	2.18	62,63,63,63	0
2	GLC	A	1093	11/12	0.89	0.19	1.72	26,30,33,35	0
2	GLC	A	1098	12/12	0.94	0.10	1.31	18,23,26,31	0
2	GLC	A	1100	12/12	0.75	0.22	0.82	60,61,62,62	0
3	CA	A	2402	1/1	0.95	0.11	0.31	35,35,35,35	0
2	GLC	A	1094	12/12	0.96	0.06	-1.06	19,22,23,23	0
3	CA	A	2405	1/1	0.84	0.14	-1.07	60,60,60,60	0
2	GLC	A	1097	11/12	0.98	0.05	-1.16	14,15,17,17	0
3	CA	A	2401	1/1	1.00	0.05	-1.47	16,16,16,16	0
3	CA	A	2403	1/1	0.99	0.06	-1.73	30,30,30,30	0
3	CA	A	2404	1/1	1.00	0.03	-3.39	18,18,18,18	0
2	GLC	A	1096	11/12	0.95	0.08	-	20,24,28,29	0
2	GLC	A	1095	11/12	0.78	0.14	-	32,35,38,40	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.