



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:55 PM UTC

PDB ID : 6J35 / pdb_00006j35
Title : Crystal structure of ligand-free of PulA-G680L mutant from *Klebsiella pneumoniae*
Authors : Saka, N.; Iwamoto, H.; Takahashi, N.; Mizutani, K.; Mikami, B.
Deposited on : 2019-01-09
Resolution : 1.84 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

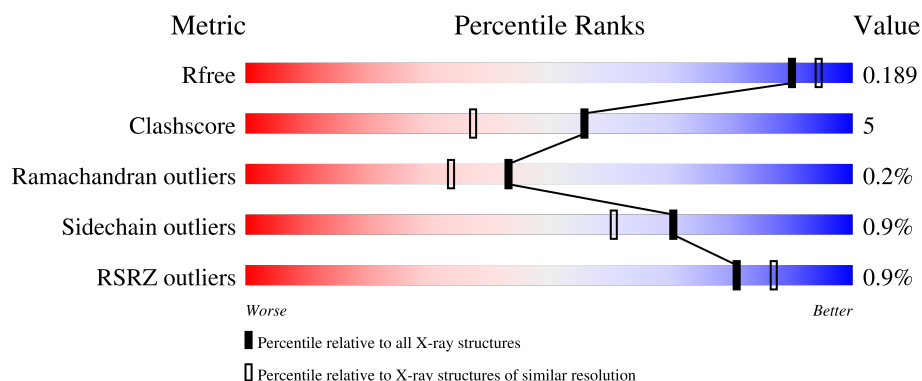
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.84 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	1296 (1.84-1.84)
Clashscore	190562	1329 (1.84-1.84)
Ramachandran outliers	187476	1318 (1.84-1.84)
Sidechain outliers	187428	1318 (1.84-1.84)
RSRZ outliers	180081	1296 (1.84-1.84)

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	1053	% 91% 8%
1	B	1053	% 91% 8% ..

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 18080 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
1	A	1052	Total	C	N	O	S	0	27	0
			8263	5145	1415	1676	27			
1	B	1045	Total	C	N	O	S	0	30	0
			8243	5136	1404	1675	28			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	31	MET	-	initiating methionine	UNP W9BQ28
A	229	SER	ASN	see sequence details	UNP W9BQ28
A	680	LEU	GLY	engineered mutation	UNP W9BQ28
B	31	MET	-	initiating methionine	UNP W9BQ28
B	229	SER	ASN	see sequence details	UNP W9BQ28
B	680	LEU	GLY	engineered mutation	UNP W9BQ28

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	5	Total	Mg	0	0
			5	5		
2	B	4	Total	Mg	0	0
			4	4		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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		

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Ca	0	0
			1	1		

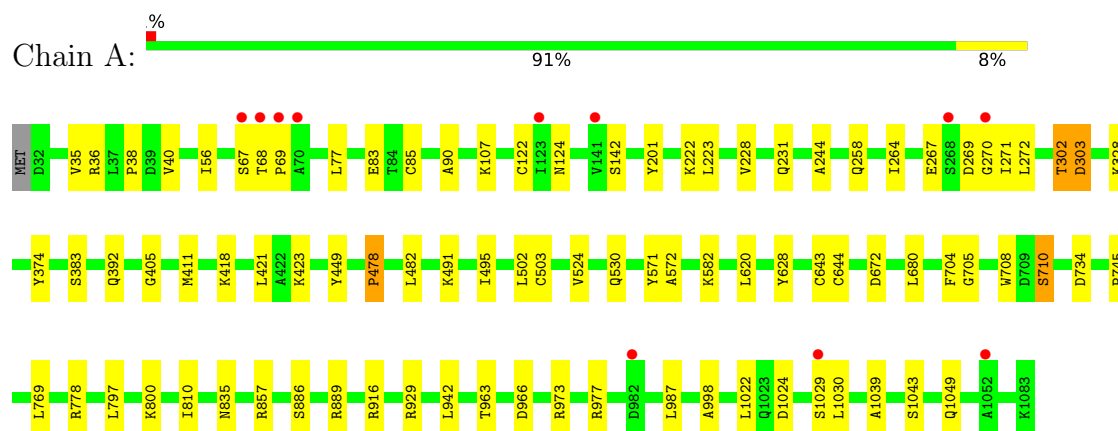
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	740	Total	O	0	0
			740	740		
5	B	812	Total	O	0	0
			812	812		

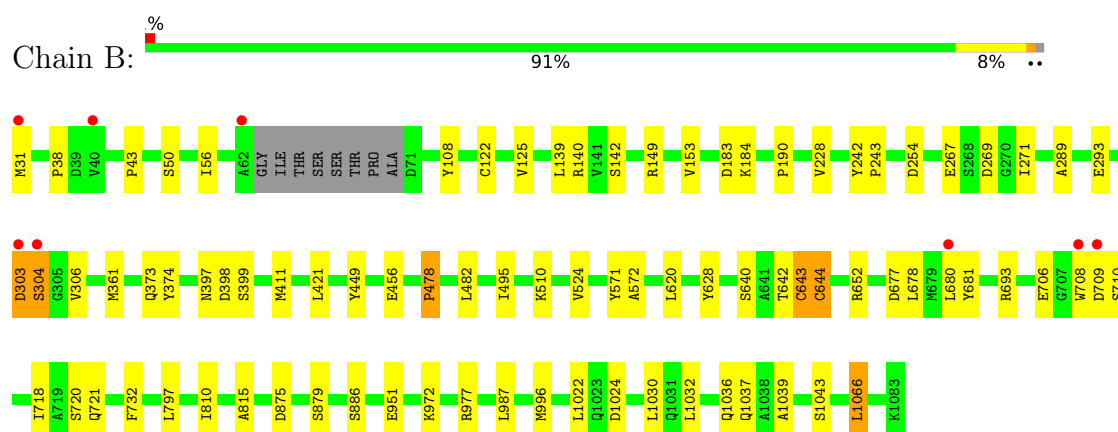
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Pullulanase



• Molecule 1: Pullulanase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	60.16Å 80.20Å 127.30Å 94.92° 103.01° 112.03°	Depositor
Resolution (Å)	38.72 – 1.84 38.72 – 1.84	Depositor EDS
% Data completeness (in resolution range)	91.1 (38.72-1.84) 91.1 (38.72-1.84)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	9.50 (at 1.84Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.164 , 0.190 0.165 , 0.189	Depositor DCC
R_{free} test set	8229 reflections (4.80%)	wwPDB-VP
Wilson B-factor (Å ²)	17.6	Xtriage
Anisotropy	1.077	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 36.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.127 for h,-h-k,-h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	18080	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.22% 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, MG, CA

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	0.54	13/8434 (0.2%)	0.60	3/11478 (0.0%)
1	B	0.53	8/8412 (0.1%)	0.61	5/11448 (0.0%)
All	All	0.54	21/16846 (0.1%)	0.60	8/22926 (0.0%)

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	90	ALA	C-O	-9.62	1.19	1.23
1	A	857	ARG	N-CA	7.41	1.55	1.46
1	B	640[A]	SER	C-O	-6.78	1.16	1.23
1	B	640[B]	SER	C-O	-6.78	1.16	1.23
1	B	304	SER	CA-C	6.44	1.61	1.52
1	A	1049	GLN	C-O	-6.16	1.16	1.24
1	A	710[A]	SER	CA-C	5.93	1.60	1.52
1	A	710[B]	SER	CA-C	5.93	1.60	1.52
1	A	503	CYS	C-O	-5.86	1.17	1.24
1	A	83	GLU	C-O	-5.77	1.16	1.24
1	A	704	PHE	C-O	-5.76	1.17	1.23
1	A	502	LEU	C-O	-5.61	1.17	1.24
1	B	306	VAL	C-O	-5.55	1.17	1.23
1	B	398	ASP	C-O	-5.51	1.17	1.23
1	A	1030	LEU	C-O	-5.33	1.17	1.23
1	B	399	SER	C-O	-5.26	1.17	1.24
1	A	705	GLY	C-O	-5.24	1.18	1.23
1	A	1029	SER	C-O	-5.12	1.16	1.24
1	B	242	TYR	C-O	-5.11	1.17	1.23
1	A	491	LYS	C-O	-5.06	1.17	1.24
1	B	397	ASN	C-O	-5.04	1.17	1.24

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	304	SER	N-CA-C	-10.54	100.16	113.23
1	B	303	ASP	N-CA-C	8.87	124.36	112.30
1	A	302	THR	CB-CA-C	-6.96	100.59	110.34
1	B	642	THR	CA-C-N	5.74	132.49	121.54
1	B	642	THR	C-N-CA	5.74	132.49	121.54
1	A	710[A]	SER	CA-C-O	5.64	129.00	121.78
1	A	710[B]	SER	CA-C-O	5.64	129.00	121.78
1	B	303	ASP	CA-C-O	5.49	126.15	119.78

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	8263	0	7923	68	0
1	B	8243	0	7879	86	0
2	A	5	0	0	0	0
2	B	4	0	0	0	0
3	A	12	0	16	3	0
4	B	1	0	0	0	0
5	A	740	0	0	18	0
5	B	812	0	0	12	0
All	All	18080	0	15818	150	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 5.

All (150) 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:680[A]:LEU:HD12	1:B:710[A]:SER:CB	1.56	1.33
1:B:680[B]:LEU:CD2	1:B:709[B]:ASP:O	1.80	1.28
1:B:228:VAL:HG23	5:B:1345:HOH:O	1.45	1.12
1:B:680[A]:LEU:CD1	1:B:710[A]:SER:HB3	1.80	1.11

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:680[B]:LEU:HD23	1:B:709[B]:ASP:O	1.40	1.07
1:B:708[B]:TRP:CZ2	1:B:720:SER:HA	2.04	0.93
1:A:411[B]:MET:HE1	5:A:1292:HOH:O	1.67	0.93
1:B:680[B]:LEU:HD21	1:B:709[B]:ASP:O	1.66	0.93
1:B:706[B]:GLU:HB3	1:B:732:PHE:CD1	2.06	0.90
1:A:328[B]:LYS:HE2	5:A:1220:HOH:O	1.73	0.88
1:B:680[A]:LEU:HD12	1:B:710[A]:SER:HB3	0.87	0.86
1:B:708[B]:TRP:HZ2	1:B:721:GLN:N	1.75	0.84
1:B:706[B]:GLU:HB3	1:B:732:PHE:HD1	1.40	0.82
1:B:708[B]:TRP:CZ2	1:B:721:GLN:N	2.49	0.80
1:B:680[A]:LEU:CD1	1:B:710[A]:SER:CB	2.51	0.79
1:A:328[B]:LYS:CE	5:A:1220:HOH:O	2.33	0.76
1:A:68:THR:HB	1:B:709[A]:ASP:HB2	1.70	0.73
1:B:678:LEU:HD23	5:B:1291:HOH:O	1.88	0.72
1:A:977:ARG:HH11	1:A:1024:ASP:HB3	1.56	0.68
1:B:680[A]:LEU:HD12	1:B:710[A]:SER:HB2	1.67	0.68
1:B:987:LEU:HD21	1:B:1022:LEU:HD21	1.74	0.68
1:A:644:CYS:SG	5:A:1594:HOH:O	2.51	0.68
1:B:708[B]:TRP:HZ2	1:B:720:SER:C	2.03	0.67
1:A:942:LEU:CD1	5:A:1642:HOH:O	2.42	0.67
1:A:987:LEU:HD21	1:A:1022:LEU:HD21	1.77	0.65
1:B:706[B]:GLU:HA	1:B:708[B]:TRP:HZ3	1.61	0.65
1:A:411[B]:MET:CE	5:A:1292:HOH:O	2.34	0.64
1:B:977:ARG:HH11	1:B:1024:ASP:HB3	1.64	0.62
1:B:708[B]:TRP:CZ2	1:B:720:SER:CA	2.80	0.62
1:B:708[B]:TRP:HZ2	1:B:720:SER:HA	1.66	0.59
1:B:289:ALA:O	1:B:293[B]:GLU:HG3	2.03	0.59
1:A:36[B]:ARG:NH1	1:A:269:ASP:O	2.36	0.58
1:B:708[B]:TRP:CE3	1:B:708[B]:TRP:N	2.72	0.58
1:B:293[B]:GLU:OE2	1:B:652:ARG:HB3	2.04	0.57
1:B:710[B]:SER:HB2	1:B:718:ILE:CD1	2.34	0.57
1:A:68:THR:CB	1:B:709[A]:ASP:HB2	2.35	0.57
1:B:706[B]:GLU:HA	1:B:708[B]:TRP:CZ3	2.40	0.56
1:A:916:ARG:NH2	5:A:1209:HOH:O	2.35	0.56
1:B:708[B]:TRP:HZ2	1:B:720:SER:CA	2.19	0.56
1:B:693[B]:ARG:NH1	5:B:1204:HOH:O	2.37	0.55
1:B:708[B]:TRP:HE3	1:B:708[B]:TRP:H	1.54	0.55
1:B:680[B]:LEU:C	1:B:680[B]:LEU:HD22	2.32	0.55
1:A:383:SER:HB3	1:A:392:GLN:HB3	1.90	0.54
1:A:582:LYS:HB2	5:A:1260:HOH:O	2.07	0.54
1:A:228:VAL:HG12	1:A:244:ALA:HB2	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1039:ALA:HB3	1:B:1043:SER:HB2	1.89	0.54
1:A:271:ILE:HD12	1:A:272:LEU:H	1.73	0.53
1:B:38:PRO:HA	1:B:271:ILE:HD12	1.90	0.53
1:B:708[B]:TRP:N	1:B:708[B]:TRP:HE3	2.05	0.53
1:A:258:GLN:NE2	5:A:1213:HOH:O	2.39	0.53
1:A:977:ARG:NH1	1:A:1024:ASP:HB3	2.21	0.53
1:A:302:THR:O	1:A:302:THR:OG1	2.21	0.53
1:A:68:THR:CG2	1:B:709[A]:ASP:HB2	2.39	0.52
1:B:680[B]:LEU:CD1	1:B:680[B]:LEU:N	2.73	0.52
1:B:706[B]:GLU:CB	1:B:732:PHE:CD1	2.88	0.52
1:A:745:PRO:HD2	5:A:1285:HOH:O	2.10	0.51
1:A:680:LEU:HG	1:A:710[B]:SER:OG	2.10	0.51
1:B:495:ILE:HA	1:B:524:VAL:HB	1.93	0.51
1:B:708[B]:TRP:NE1	1:B:721:GLN:H	2.07	0.51
1:B:680[B]:LEU:CD1	1:B:680[B]:LEU:H	2.23	0.51
1:B:680[B]:LEU:HD13	1:B:681[B]:TYR:H	1.75	0.51
1:B:709[B]:ASP:N	1:B:709[B]:ASP:OD1	2.43	0.51
1:B:797:LEU:HD11	1:B:810:ILE:HD11	1.91	0.51
1:A:69:PRO:HB3	5:B:1894:HOH:O	2.11	0.51
1:A:223[B]:LEU:N	1:A:223[B]:LEU:HD13	2.26	0.51
1:B:680[B]:LEU:CG	1:B:709[B]:ASP:O	2.56	0.51
1:A:495:ILE:HA	1:A:524:VAL:HB	1.93	0.51
1:A:1039:ALA:HB3	1:A:1043:SER:HB2	1.92	0.50
1:A:222:LYS:C	1:A:223[B]:LEU:HD13	2.37	0.50
1:A:267:GLU:HB2	1:A:271:ILE:HG23	1.94	0.50
1:A:680:LEU:CD1	1:A:710[B]:SER:OG	2.60	0.50
1:A:68:THR:CG2	1:B:709[A]:ASP:CB	2.89	0.50
1:B:267:GLU:HB3	5:B:1201:HOH:O	2.11	0.49
1:A:269:ASP:OD1	1:A:271:ILE:HG22	2.12	0.49
1:B:620:LEU:HB3	1:B:628:TYR:CD2	2.48	0.49
1:B:977:ARG:NH1	1:B:1024:ASP:HB3	2.28	0.49
1:A:797:LEU:HD11	1:A:810:ILE:HD11	1.94	0.49
1:B:43:PRO:HB3	1:B:108:TYR:CD1	2.47	0.48
1:B:681[B]:TYR:OH	1:B:709[B]:ASP:HB2	2.13	0.48
1:B:951:GLU:OE1	1:B:1036:GLN:HG2	2.14	0.48
1:A:421:LEU:HD21	5:A:1477:HOH:O	2.12	0.48
1:A:269:ASP:OD1	1:A:269:ASP:N	2.47	0.48
1:B:456:GLU:OE1	1:B:456:GLU:N	2.35	0.48
1:B:680[B]:LEU:N	1:B:680[B]:LEU:HD13	2.28	0.48
1:A:418:LYS:HE2	5:A:1249:HOH:O	2.13	0.48
1:A:271:ILE:HD12	1:A:272:LEU:N	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:643:CYS:SG	1:A:644:CYS:N	2.87	0.47
1:A:644:CYS:HB3	5:A:1607:HOH:O	2.14	0.47
1:A:734:ASP:OD1	1:A:734:ASP:N	2.44	0.47
1:B:693[A]:ARG:CZ	5:B:1301:HOH:O	2.61	0.47
1:B:680[B]:LEU:HD13	1:B:681[B]:TYR:N	2.30	0.47
1:A:67:SER:O	1:B:815:ALA:HB2	2.15	0.46
1:A:680:LEU:HD13	1:A:708:TRP:HB2	1.96	0.46
1:B:1030:LEU:HG	1:B:1066:LEU:HB3	1.97	0.46
1:B:50:SER:O	1:B:149:ARG:NH1	2.48	0.46
1:B:706[B]:GLU:HB2	1:B:732:PHE:HB2	1.97	0.46
1:A:405:GLY:H	3:A:1107:GOL:C2	2.28	0.46
1:A:411[A]:MET:HE2	1:A:672:ASP:OD1	2.15	0.46
1:A:328[B]:LYS:NZ	5:A:1220:HOH:O	2.43	0.46
1:A:418:LYS:HD2	1:A:966[B]:ASP:OD1	2.16	0.46
3:A:1106:GOL:H12	5:A:1542:HOH:O	2.17	0.45
1:A:778[B]:ARG:NH1	1:A:800:LYS:O	2.48	0.45
1:B:190:PRO:HA	1:B:243:PRO:HG2	1.98	0.45
1:B:269:ASP:OD1	1:B:271:ILE:HG12	2.16	0.45
1:A:69:PRO:HB3	5:B:1846:HOH:O	2.16	0.45
1:B:708[B]:TRP:HE1	1:B:720:SER:HB2	1.81	0.45
1:B:411:MET:HE1	5:B:1575:HOH:O	2.17	0.44
1:A:680:LEU:HD12	1:A:710[B]:SER:OG	2.16	0.44
1:B:1032:LEU:HD23	1:B:1037[B]:GLN:HG2	1.97	0.44
1:A:38:PRO:HB2	1:A:40:VAL:HG12	1.99	0.44
1:A:85:CYS:HB2	1:A:124:ASN:ND2	2.33	0.44
1:A:572:ALA:HB1	5:A:1556:HOH:O	2.16	0.44
1:B:680[A]:LEU:H	1:B:680[A]:LEU:HD23	1.82	0.44
1:A:405:GLY:CA	3:A:1107:GOL:H31	2.49	0.43
1:A:973:ARG:HG2	1:A:998:ALA:HB1	2.00	0.43
1:B:643:CYS:SG	1:B:644:CYS:N	2.91	0.43
1:A:223[B]:LEU:N	1:A:223[B]:LEU:CD1	2.80	0.43
1:B:139:LEU:HD13	1:B:153:VAL:HG11	2.01	0.43
1:B:875:ASP:OD2	1:B:879:SER:HB2	2.18	0.43
1:B:421:LEU:HD11	5:B:2007:HOH:O	2.18	0.43
1:A:835:ASN:HA	1:A:889[A]:ARG:HG2	1.99	0.43
1:A:231[A]:GLN:HG2	5:A:1787:HOH:O	2.18	0.43
1:B:510:LYS:HE2	5:B:1950:HOH:O	2.17	0.43
1:A:201:TYR:CD1	1:A:264:ILE:HD12	2.54	0.43
1:A:40:VAL:HG21	1:A:107:LYS:HD2	2.01	0.42
1:A:35:VAL:HA	1:A:270:GLY:O	2.20	0.42
1:B:972:LYS:HE2	1:B:972:LYS:HB3	1.81	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:449:TYR:CZ	1:A:571:TYR:HB2	2.54	0.42
1:B:449:TYR:CZ	1:B:571:TYR:HB2	2.54	0.42
1:B:996:MET:HE3	1:B:996:MET:HB2	1.85	0.42
1:A:769:LEU:HD11	1:A:929:ARG:NH2	2.34	0.42
1:B:680[B]:LEU:H	1:B:680[B]:LEU:HD12	1.85	0.42
1:B:708[B]:TRP:NE1	1:B:720:SER:HB2	2.35	0.42
1:B:678:LEU:HB3	1:B:681[B]:TYR:HD2	1.85	0.41
1:A:267:GLU:HB2	1:A:271:ILE:CG2	2.50	0.41
1:B:43:PRO:HB3	1:B:108:TYR:CG	2.55	0.41
1:B:56:ILE:HD11	1:B:125:VAL:HG21	2.02	0.41
1:B:122:CYS:HB3	1:B:142:SER:HA	2.02	0.41
1:B:680[B]:LEU:HD23	1:B:710[B]:SER:HA	2.02	0.41
1:A:620:LEU:HB3	1:A:628:TYR:CD2	2.55	0.41
1:A:530:GLN:HB2	5:A:1653:HOH:O	2.20	0.41
1:B:361:MET:O	1:B:373:GLN:HA	2.21	0.41
1:A:222:LYS:C	1:A:223[B]:LEU:CD1	2.94	0.41
1:B:677:ASP:O	1:B:678:LEU:C	2.64	0.41
1:A:122:CYS:HB3	1:A:142:SER:HA	2.02	0.40
1:B:183:ASP:O	1:B:184:LYS:C	2.65	0.40
1:B:254:ASP:HB2	5:B:1804:HOH:O	2.19	0.40
1:B:572:ALA:HB1	5:B:1718:HOH:O	2.22	0.40
1:A:56:ILE:HG21	1:A:77:LEU:HD11	2.03	0.40
1:A:423:LYS:HD2	1:A:963:THR:OG1	2.22	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	1077/1053 (102%)	1053 (98%)	22 (2%)	2 (0%)	43	34
1	B	1071/1053 (102%)	1048 (98%)	21 (2%)	2 (0%)	43	34

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	2148/2106 (102%)	2101 (98%)	43 (2%)	4 (0%)	43	34

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	303	ASP
1	B	643	CYS
1	B	478	PRO
1	A	478	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	893/867 (103%)	888 (99%)	5 (1%)	78	72
1	B	890/867 (103%)	880 (99%)	10 (1%)	65	54
All	All	1783/1734 (103%)	1768 (99%)	15 (1%)	70	65

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	303	ASP
1	A	374	TYR
1	A	478	PRO
1	A	482	LEU
1	A	886	SER
1	B	31	MET
1	B	140	ARG
1	B	303	ASP
1	B	304	SER
1	B	374	TYR
1	B	478	PRO
1	B	482	LEU
1	B	644	CYS
1	B	886	SER

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Mol	Chain	Res	Type
1	B	1066	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	230	GLN
1	A	465	GLN
1	A	593	GLN
1	B	96	ASN
1	B	124	ASN
1	B	194	ASN
1	B	373	GLN
1	B	439	GLN
1	B	461	GLN
1	B	1049	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](#)

Of 12 ligands modelled in this entry, 10 are monoatomic - leaving 2 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
3	GOL	A	1107	-	5,5,5	0.45	0	5,5,5	0.84	0
3	GOL	A	1106	-	5,5,5	0.50	0	5,5,5	0.90	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	1107	-	-	2/4/4/4	-
3	GOL	A	1106	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1106	GOL	C1-C2-C3-O3
3	A	1107	GOL	C1-C2-C3-O3
3	A	1106	GOL	O2-C2-C3-O3
3	A	1107	GOL	O2-C2-C3-O3

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1107	GOL	2	0
3	A	1106	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	1052/1053 (99%)	-0.10	11 (1%) 79 86	8, 19, 39, 55	27 (2%)
1	B	1045/1053 (99%)	-0.19	8 (0%) 82 88	8, 18, 35, 62	30 (2%)
All	All	2097/2106 (99%)	-0.14	19 (0%) 81 87	8, 19, 38, 62	57 (2%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	31	MET	4.8
1	B	708[A]	TRP	3.8
1	B	62	ALA	3.7
1	B	709[A]	ASP	3.3
1	A	68	THR	3.3
1	A	70	ALA	2.6
1	A	67	SER	2.5
1	B	680[A]	LEU	2.4
1	A	268	SER	2.3
1	A	69	PRO	2.3
1	B	40	VAL	2.3
1	B	304	SER	2.3
1	A	270	GLY	2.2
1	A	1029	SER	2.2
1	A	1052	ALA	2.2
1	A	982	ASP	2.1
1	B	303	ASP	2.1
1	A	123	ILE	2.0
1	A	141	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 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
3	GOL	A	1107	6/6	0.77	0.15	20,20,20,20	6
3	GOL	A	1106	6/6	0.81	0.19	20,20,20,20	6
2	MG	A	1105	1/1	0.92	0.13	32,32,32,32	0
2	MG	A	1104	1/1	0.97	0.04	28,28,28,28	0
2	MG	B	1103	1/1	0.97	0.07	26,26,26,26	0
2	MG	B	1102	1/1	0.98	0.03	18,18,18,18	0
2	MG	B	1104	1/1	0.98	0.05	26,26,26,26	0
2	MG	A	1101	1/1	0.99	0.11	11,11,11,11	0
2	MG	A	1102	1/1	0.99	0.10	11,11,11,11	0
2	MG	B	1101	1/1	0.99	0.09	10,10,10,10	0
2	MG	A	1103	1/1	0.99	0.02	20,20,20,20	0
4	CA	B	1105	1/1	0.99	0.09	30,30,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.