



Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 01:30 am BST

PDB ID : 6J33
Title : Crystal structure of ligand-free of PulA from *Klebsiella pneumoniae*
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Deposited on : 2019-01-09
Resolution : 1.30 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

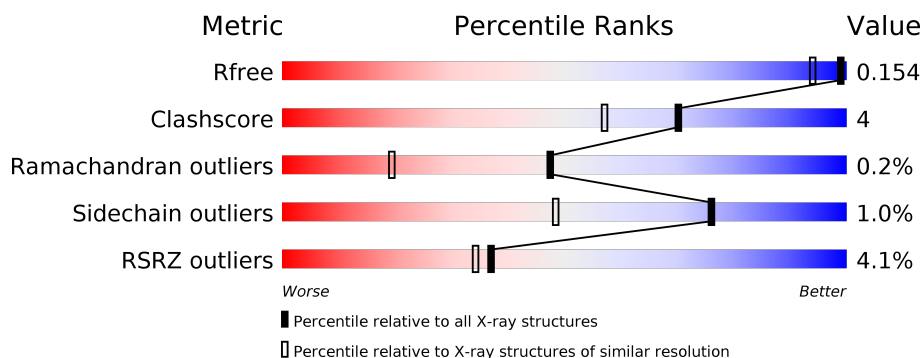
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.30 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	1058 (1.30-1.30)
Clashscore	141614	1101 (1.30-1.30)
Ramachandran outliers	138981	1058 (1.30-1.30)
Sidechain outliers	138945	1058 (1.30-1.30)
RSRZ outliers	127900	1029 (1.30-1.30)

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	1053	<div> <div>5%</div> <div> <div></div> <div>94%</div> <div>6%</div> </div> </div>
1	B	1053	<div> <div>3%</div> <div> <div></div> <div>94%</div> <div>6%</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 20463 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	1045	Total	C	N	O	S	0	70	0
			8553	5313	1469	1744	27			
1	B	1053	Total	C	N	O	S	0	84	0
			8723	5409	1500	1787	27			

There are 4 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
B	31	MET	-	initiating methionine	UNP W9BQ28
B	229	SER	ASN	see sequence details	UNP W9BQ28

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	10	Total	Mg	0	0
			10	10		
2	A	12	Total	Mg	0	0
			12	12		

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

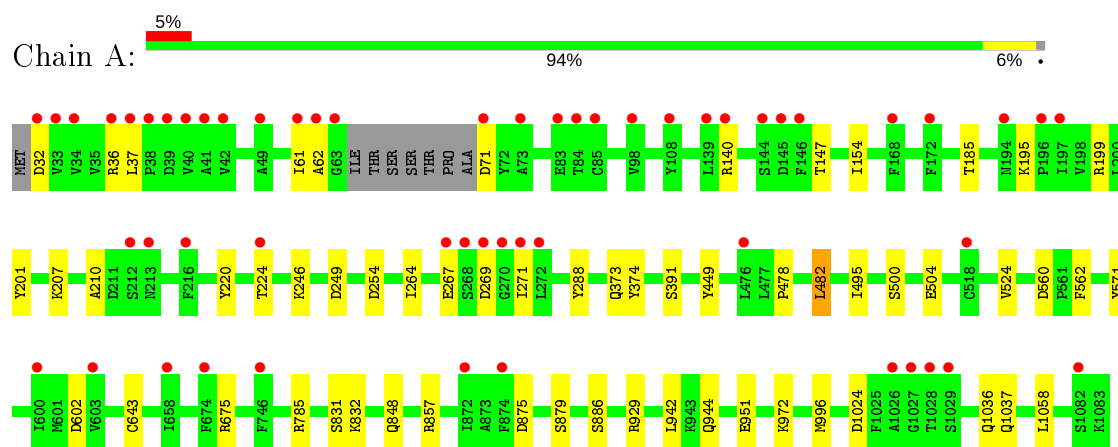
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1527	Total 1527	O 1527	0	0
5	B	1605	Total 1605	O 1605	0	0

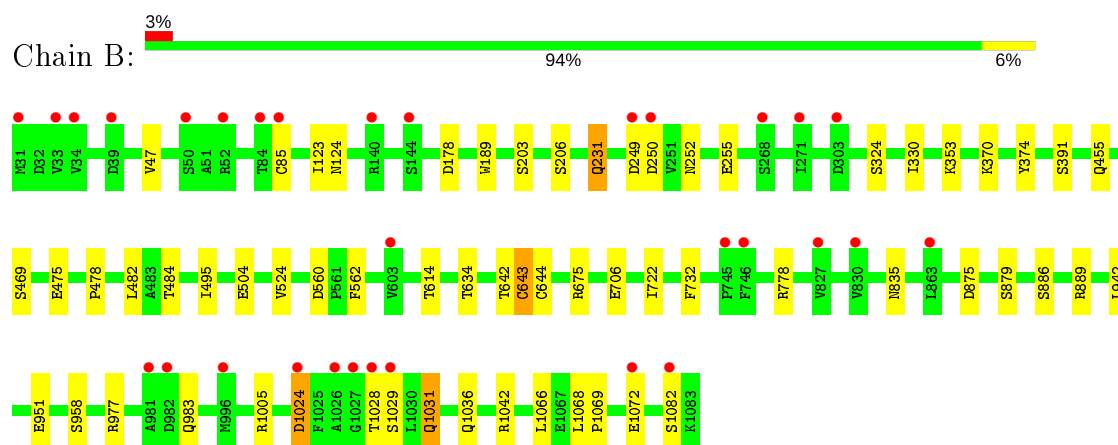
3 Residue-property plots [i](#)

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: 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	60.36Å 80.28Å 127.74Å 96.70° 102.30° 112.10°	Depositor
Resolution (Å)	39.26 – 1.30 39.26 – 1.30	Depositor EDS
% Data completeness (in resolution range)	94.4 (39.26-1.30) 94.4 (39.26-1.30)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.49 (at 1.30Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.128 , 0.154 0.128 , 0.154	Depositor DCC
R_{free} test set	24855 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	15.0	Xtriage
Anisotropy	0.173	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 55.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.064 for h,-h-k,-h-l	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	20463	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.26% 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

5.1 Standard geometry

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

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	0.43	0/8723	0.63	0/11856
1	B	0.39	0/8897	0.61	0/12097
All	All	0.41	0/17620	0.62	0/23953

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	8553	0	8168	52	0
1	B	8723	0	8304	66	0
2	A	12	0	0	0	0
2	B	10	0	0	0	0
3	A	12	0	9	1	0
3	B	20	0	15	0	0
4	B	1	0	0	0	0
5	A	1527	0	0	31	0
5	B	1605	0	0	28	0
All	All	20463	0	16496	119	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 4.

All (119) 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)
3:A:1113:ACT:H2	5:A:1794:HOH:O	1.47	1.13
1:A:36[B]:ARG:HH21	1:A:271:ILE:HG22	1.25	1.01
1:B:504[A]:GLU:HG2	5:B:1555:HOH:O	1.62	1.00
1:A:560:ASP:HB2	5:A:1888:HOH:O	1.64	0.95
1:B:1082[B]:SER:HB2	5:B:1268:HOH:O	1.70	0.91
1:B:614[A]:THR:HG22	5:B:1813:HOH:O	1.71	0.90
1:A:36[B]:ARG:NH2	1:A:271:ILE:HG22	1.91	0.84
1:B:977:ARG:NH1	1:B:1024[B]:ASP:HB3	1.95	0.82
1:A:942[A]:LEU:HD21	5:A:1995:HOH:O	1.84	0.78
1:B:675:ARG:NH2	1:B:706:GLU:HG3	1.98	0.78
1:A:482[A]:LEU:HB2	5:A:1888:HOH:O	1.83	0.77
1:A:373[B]:GLN:HG3	5:A:2120:HOH:O	1.88	0.73
1:B:942[A]:LEU:HD21	5:B:1882:HOH:O	1.88	0.73
1:A:972[A]:LYS:HD2	5:A:2281:HOH:O	1.89	0.73
1:B:977:ARG:HH11	1:B:1024[B]:ASP:HB3	1.55	0.72
1:A:254[B]:ASP:HB2	5:A:1558:HOH:O	1.88	0.71
1:B:675:ARG:HD2	1:B:732:PHE:HE2	1.56	0.71
1:A:785[B]:ARG:NH1	5:A:1201:HOH:O	2.23	0.70
1:A:942[A]:LEU:CD2	5:A:1995:HOH:O	2.38	0.70
1:A:675:ARG:HH21	1:A:832[B]:LYS:NZ	1.90	0.70
1:B:1031[B]:GLN:OE1	1:B:1069:PRO:CG	2.40	0.69
1:A:1058:LEU:HD22	5:A:2307:HOH:O	1.92	0.69
1:B:675:ARG:HD2	1:B:732:PHE:CE2	2.28	0.68
1:B:1031[B]:GLN:OE1	1:B:1069:PRO:HG2	1.95	0.67
1:A:972[A]:LYS:CD	5:A:2281:HOH:O	2.43	0.65
1:A:269:ASP:OD1	1:A:271:ILE:HG23	1.96	0.65
1:A:1058:LEU:HD13	5:A:2307:HOH:O	1.95	0.65
1:A:848:GLN:NE2	5:A:1205:HOH:O	2.30	0.65
1:A:1037[B]:GLN:NE2	5:A:1204:HOH:O	2.30	0.64
1:B:1082[A]:SER:HA	5:B:2107:HOH:O	1.95	0.64
1:B:942[A]:LEU:CD2	5:B:1882:HOH:O	2.44	0.63
1:A:929[A]:ARG:HD2	5:A:2229:HOH:O	1.98	0.62
1:A:972[A]:LYS:CG	5:A:2281:HOH:O	2.46	0.62
1:B:47[A]:VAL:HG23	5:B:2423:HOH:O	1.99	0.61
1:A:207[A]:LYS:HE3	5:A:1311:HOH:O	2.00	0.61
1:B:977:ARG:NH1	1:B:1024[A]:ASP:HB2	2.16	0.61
1:B:370[B]:LYS:HE3	5:B:2133:HOH:O	2.00	0.61
1:B:1031[B]:GLN:OE1	1:B:1069:PRO:HG3	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1058:LEU:HB3	5:A:2307:HOH:O	2.02	0.60
1:B:1028:THR:HG22	1:B:1068:LEU:HD21	1.84	0.60
1:B:675:ARG:HH22	1:B:706:GLU:HG3	1.64	0.59
1:A:944:GLN:NE2	5:A:1207:HOH:O	2.34	0.59
1:B:958:SER:O	1:B:1005[A]:ARG:NH1	2.31	0.59
1:B:634[B]:THR:HG21	5:B:1837:HOH:O	2.02	0.58
1:A:951[A]:GLU:HG3	5:A:1648:HOH:O	2.04	0.58
1:B:644:CYS:HB3	5:B:2049:HOH:O	2.04	0.58
1:B:504[B]:GLU:HG3	5:B:1555:HOH:O	2.04	0.57
1:A:675:ARG:HH21	1:A:832[B]:LYS:HZ3	1.53	0.57
1:A:147:THR:HG23	5:A:1216:HOH:O	2.04	0.57
1:A:972[A]:LYS:HG3	5:A:2281:HOH:O	2.05	0.57
1:B:1082[B]:SER:HA	5:B:2107:HOH:O	2.03	0.57
1:A:195:LYS:HE3	1:A:267:GLU:OE2	2.05	0.56
1:A:36[B]:ARG:NH2	1:A:271:ILE:CG2	2.67	0.56
1:B:835:ASN:ND2	5:B:1215:HOH:O	2.40	0.55
1:A:246[A]:LYS:NZ	5:A:1214:HOH:O	2.40	0.55
1:B:643:CYS:SG	1:B:644:CYS:N	2.81	0.53
1:B:889[A]:ARG:NH2	5:B:1203:HOH:O	2.30	0.53
1:B:1005[A]:ARG:CZ	5:B:1283:HOH:O	2.57	0.52
1:B:675:ARG:HH22	1:B:706:GLU:CG	2.22	0.52
1:B:675:ARG:HH22	1:B:706:GLU:CD	2.13	0.52
1:A:61:ILE:HG13	1:A:154:ILE:HD13	1.91	0.51
1:A:929[A]:ARG:CD	5:A:2229:HOH:O	2.57	0.51
1:B:778[A]:ARG:NH2	1:B:983:GLN:O	2.43	0.51
1:A:875:ASP:OD2	1:A:879:SER:HB2	2.11	0.50
1:B:1042:ARG:HD2	5:B:1273:HOH:O	2.10	0.50
1:B:203[B]:SER:OG	1:B:206[B]:SER:O	2.20	0.50
1:B:1028:THR:CG2	1:B:1068:LEU:HD21	2.40	0.50
1:A:254[B]:ASP:OD2	1:A:288:TYR:OH	2.29	0.49
1:B:889[A]:ARG:NH1	5:B:1238:HOH:O	2.44	0.49
1:A:224[B]:THR:HG23	5:A:1813:HOH:O	2.12	0.49
1:A:36[B]:ARG:HG2	1:A:37:LEU:O	2.12	0.49
1:A:37:LEU:HD21	1:A:210:ALA:CB	2.42	0.49
1:B:1024[A]:ASP:N	1:B:1024[A]:ASP:OD1	2.39	0.49
1:B:469[A]:SER:HB2	5:B:1359:HOH:O	2.13	0.49
1:B:484:THR:OG1	1:B:560[B]:ASP:OD2	2.28	0.48
1:B:951[A]:GLU:OE2	1:B:1036:GLN:HG2	2.13	0.48
1:B:875:ASP:OD2	1:B:879:SER:HB2	2.14	0.48
1:B:495:ILE:HA	1:B:524:VAL:HB	1.96	0.48
1:A:482[B]:LEU:HB3	5:A:1888:HOH:O	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:231[A]:GLN:H	1:B:231[A]:GLN:CD	2.17	0.47
1:B:1028:THR:HG22	1:B:1068:LEU:CD2	2.45	0.47
1:B:722:ILE:HG13	5:B:2213:HOH:O	2.14	0.47
1:A:495:ILE:HA	1:A:524:VAL:HB	1.97	0.47
1:A:207[B]:LYS:HE2	5:A:1655:HOH:O	2.15	0.47
1:B:1028:THR:O	1:B:1028:THR:HG22	2.14	0.47
1:A:199:ARG:HD3	1:A:220:TYR:CD1	2.50	0.46
1:B:252:ASN:O	1:B:255:GLU:HB2	2.15	0.46
1:B:889[A]:ARG:NH1	5:B:1203:HOH:O	2.40	0.46
1:B:324:SER:HB3	1:B:330[B]:ILE:HD11	1.97	0.46
1:B:675:ARG:CZ	5:B:1938:HOH:O	2.64	0.45
1:A:32:ASP:N	5:A:1242:HOH:O	2.48	0.45
1:B:475:GLU:OE2	1:B:675:ARG:HD3	2.16	0.45
1:A:951[A]:GLU:OE2	1:A:1036:GLN:HG2	2.16	0.45
1:B:123:ILE:HA	5:B:2189:HOH:O	2.16	0.45
1:B:353:LYS:NZ	5:B:1250:HOH:O	2.48	0.45
1:A:996:MET:HE3	5:A:2626:HOH:O	2.16	0.44
1:B:1082[A]:SER:CB	5:B:1268:HOH:O	2.65	0.44
1:B:675:ARG:NH2	1:B:706:GLU:CG	2.76	0.43
1:A:602[A]:ASP:OD1	1:A:675:ARG:HD2	2.18	0.43
1:B:1082[A]:SER:HB3	5:B:1268:HOH:O	2.19	0.43
1:B:85:CYS:HB2	1:B:124:ASN:ND2	2.34	0.43
1:B:1005[A]:ARG:HB2	1:B:1005[A]:ARG:HE	1.66	0.42
1:A:1024:ASP:N	1:A:1024:ASP:OD1	2.48	0.42
1:A:500:SER:O	1:A:504[A]:GLU:HG3	2.19	0.42
1:A:185:THR:O	1:A:246[B]:LYS:HA	2.20	0.42
1:B:178[B]:ASP:OD1	1:B:189:TRP:NE1	2.51	0.42
1:B:250[A]:ASP:OD1	1:B:250[A]:ASP:N	2.53	0.42
1:A:246[B]:LYS:NZ	1:A:249:ASP:OD1	2.53	0.41
1:A:449:TYR:CZ	1:A:571:TYR:HB2	2.55	0.41
1:A:857[B]:ARG:NH1	5:A:1254:HOH:O	2.51	0.41
1:B:675:ARG:HH21	1:B:706:GLU:HG3	1.81	0.41
1:B:1028:THR:O	1:B:1028:THR:CG2	2.69	0.40
1:A:140[B]:ARG:NH1	5:A:1217:HOH:O	2.40	0.40
1:B:642:THR:O	1:B:643:CYS:HB3	2.22	0.40
1:B:1005[A]:ARG:NE	5:B:1283:HOH:O	2.54	0.40
1:B:1072[B]:GLU:HG3	1:B:1072[B]:GLU:H	1.72	0.40
1:B:560[B]:ASP:HA	5:B:1910:HOH:O	2.21	0.40
1:A:201:TYR:CD1	1:A:264:ILE:HD12	2.56	0.40
1:B:455:GLN:HG2	5:B:2653:HOH:O	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	1111/1053 (106%)	1092 (98%)	16 (1%)	3 (0%)	41	17
1	B	1136/1053 (108%)	1111 (98%)	23 (2%)	2 (0%)	47	19
All	All	2247/2106 (107%)	2203 (98%)	39 (2%)	5 (0%)	47	19

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	643	CYS
1	B	643	CYS
1	A	62	ALA
1	A	478	PRO
1	B	478	PRO

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	928/866 (107%)	919 (99%)	9 (1%)	76	48
1	B	950/866 (110%)	933 (98%)	17 (2%)	59	24
All	All	1878/1732 (108%)	1852 (99%)	26 (1%)	76	34

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	71	ASP

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Mol	Chain	Res	Type
1	A	374	TYR
1	A	391[A]	SER
1	A	391[B]	SER
1	A	482[A]	LEU
1	A	482[B]	LEU
1	A	562	PHE
1	A	831	SER
1	A	886	SER
1	B	231[A]	GLN
1	B	231[B]	GLN
1	B	249[A]	ASP
1	B	249[B]	ASP
1	B	374	TYR
1	B	391[A]	SER
1	B	391[B]	SER
1	B	482[A]	LEU
1	B	482[B]	LEU
1	B	562	PHE
1	B	886	SER
1	B	1024[A]	ASP
1	B	1024[B]	ASP
1	B	1029	SER
1	B	1031[A]	GLN
1	B	1031[B]	GLN
1	B	1066	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	944	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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 31 ligands modelled in this entry, 23 are monoatomic - leaving 8 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	ACT	A	1112	-	1,3,3	0.05	0	0,3,3	0.00	-
3	ACT	B	1114	-	1,3,3	1.65	0	0,3,3	0.00	-
3	ACT	B	1113	-	1,3,3	0.53	0	0,3,3	0.00	-
3	ACT	A	1113	-	1,3,3	0.40	0	0,3,3	0.00	-
3	ACT	B	1116	-	1,3,3	0.18	0	0,3,3	0.00	-
3	ACT	A	1114	-	1,3,3	0.78	0	0,3,3	0.00	-
3	ACT	B	1115	-	1,3,3	1.28	0	0,3,3	0.00	-
3	ACT	B	1112	-	1,3,3	1.25	0	0,3,3	0.00	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1113	ACT	1	0

5.7 Other polymers [i](#)

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	1045/1053 (99%)	0.25	55 (5%) 26 24	12, 17, 41, 81	1 (0%)
1	B	1053/1053 (100%)	0.10	31 (2%) 51 49	12, 17, 34, 65	0
All	All	2098/2106 (99%)	0.18	86 (4%) 37 34	12, 17, 38, 81	1 (0%)

All (86) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	40	VAL	8.5
1	A	39	ASP	8.4
1	A	62	ALA	7.6
1	A	271	ILE	7.0
1	B	746[A]	PHE	6.7
1	B	31	MET	6.6
1	A	63	GLY	6.5
1	B	1027	GLY	6.4
1	A	85	CYS	6.0
1	A	213	ASN	5.3
1	A	36[A]	ARG	5.2
1	B	249[A]	ASP	5.2
1	A	37	LEU	5.0
1	A	38	PRO	5.0
1	A	269	ASP	4.7
1	A	268	SER	4.6
1	B	1026	ALA	4.4
1	A	61	ILE	4.3
1	A	33	VAL	4.3
1	A	34	VAL	4.3
1	A	41	ALA	4.3
1	A	1028[A]	THR	4.2
1	A	212	SER	4.2
1	A	168	PHE	4.1

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Mol	Chain	Res	Type	RSRZ
1	A	42	VAL	4.0
1	B	140[A]	ARG	3.8
1	A	272	LEU	3.8
1	A	1027[A]	GLY	3.7
1	A	73	ALA	3.4
1	A	71	ASP	3.3
1	B	39[A]	ASP	3.2
1	B	85	CYS	3.2
1	A	84	THR	3.1
1	A	1026	ALA	3.1
1	B	1028	THR	3.0
1	A	270	GLY	3.0
1	A	146	PHE	3.0
1	A	197	ILE	3.0
1	B	271	ILE	3.0
1	A	108	TYR	2.9
1	A	196	PRO	2.9
1	A	476	LEU	2.7
1	B	268[A]	SER	2.7
1	B	1072[A]	GLU	2.7
1	A	746[A]	PHE	2.7
1	B	84	THR	2.6
1	A	32	ASP	2.6
1	B	1029	SER	2.6
1	B	1082[A]	SER	2.5
1	A	139	LEU	2.5
1	B	745	PRO	2.5
1	B	250[A]	ASP	2.5
1	B	830	VAL	2.5
1	A	140[A]	ARG	2.5
1	A	144	SER	2.5
1	A	216	PHE	2.5
1	B	863	LEU	2.4
1	B	827	VAL	2.4
1	A	267	GLU	2.4
1	A	1029[A]	SER	2.4
1	B	982[A]	ASP	2.4
1	A	872	ILE	2.3
1	A	98	VAL	2.3
1	B	144	SER	2.3
1	B	603	VAL	2.3
1	A	674	PHE	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	600	ILE	2.3
1	A	49	ALA	2.3
1	A	603	VAL	2.3
1	A	172	PHE	2.2
1	B	52	ARG	2.2
1	A	658	ILE	2.2
1	A	518	CYS	2.2
1	A	83[A]	GLU	2.2
1	B	1024[A]	ASP	2.2
1	A	194	ASN	2.1
1	B	303	ASP	2.1
1	B	34	VAL	2.1
1	A	145	ASP	2.1
1	B	981	ALA	2.1
1	A	874	PHE	2.1
1	A	1082	SER	2.1
1	B	996	MET	2.0
1	A	224[A]	THR	2.0
1	B	50	SER	2.0
1	B	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. 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(\AA^2)	Q<0.9
3	ACT	B	1116	4/4	0.81	0.17	56,56,56,57	0
3	ACT	B	1115	4/4	0.83	0.15	36,39,39,40	0
3	ACT	A	1114	4/4	0.85	0.23	39,41,42,44	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	ACT	A	1113	4/4	0.88	0.16	19,20,21,23	4
3	ACT	B	1112	4/4	0.90	0.13	19,21,22,25	4
3	ACT	A	1112	4/4	0.91	0.10	20,24,26,26	4
2	MG	B	1105	1/1	0.92	0.11	50,50,50,50	0
2	MG	A	1108	1/1	0.92	0.18	36,36,36,36	1
2	MG	B	1107	1/1	0.93	0.13	30,30,30,30	1
3	ACT	B	1113	4/4	0.94	0.08	21,22,22,22	4
2	MG	A	1110	1/1	0.94	0.14	26,26,26,26	1
3	ACT	B	1114	4/4	0.95	0.08	22,23,24,24	4
2	MG	A	1104	1/1	0.95	0.10	38,38,38,38	0
2	MG	A	1109	1/1	0.96	0.09	35,35,35,35	1
2	MG	B	1109	1/1	0.96	0.10	32,32,32,32	1
2	MG	A	1107	1/1	0.96	0.12	38,38,38,38	1
2	MG	B	1110	1/1	0.97	0.15	38,38,38,38	1
2	MG	B	1108	1/1	0.97	0.09	28,28,28,28	1
2	MG	B	1106	1/1	0.97	0.08	32,32,32,32	1
4	CA	B	1111	1/1	0.98	0.07	18,18,18,18	0
2	MG	A	1103	1/1	0.99	0.12	29,29,29,29	0
2	MG	B	1102	1/1	0.99	0.06	22,22,22,22	0
2	MG	A	1111	1/1	0.99	0.16	28,28,28,28	0
2	MG	B	1101	1/1	1.00	0.09	10,10,10,10	0
2	MG	A	1102	1/1	1.00	0.04	18,18,18,18	0
2	MG	B	1103	1/1	1.00	0.06	25,25,25,25	0
2	MG	A	1106	1/1	1.00	0.11	11,11,11,11	0
2	MG	A	1115	1/1	1.00	0.08	19,19,19,19	0
2	MG	A	1101	1/1	1.00	0.06	20,20,20,20	0
2	MG	B	1104	1/1	1.00	0.05	17,17,17,17	0
2	MG	A	1105	1/1	1.00	0.14	11,11,11,11	0

6.5 Other polymers [i](#)

There are no such residues in this entry.