



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:26 PM UTC

PDB ID : 6RJ1 / pdb_00006rj1
Title : N-Domain P40/P90 Mycoplasma pneumoniae
Authors : Vizarraga, D.; Aparicio, D.; Illanes, R.; Fita, I.
Deposited on : 2019-04-25
Resolution : 2.65 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

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
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

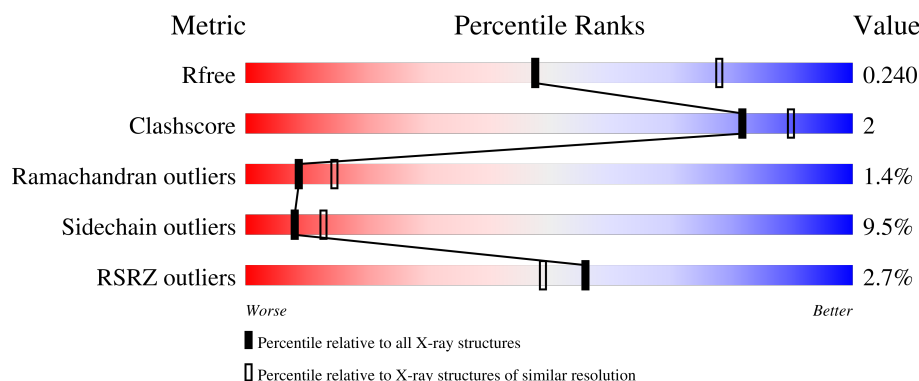
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 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}	180053	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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 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	976	2% 71% 11% • 17%
1	B	976	3% 70% 11% • 17%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12623 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 Mgp-operon protein 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	812	Total	C	N	O	S	0	2	0
			6292	3953	1076	1259	4			
1	B	806	Total	C	N	O	S	0	4	0
			6267	3944	1068	1251	4			

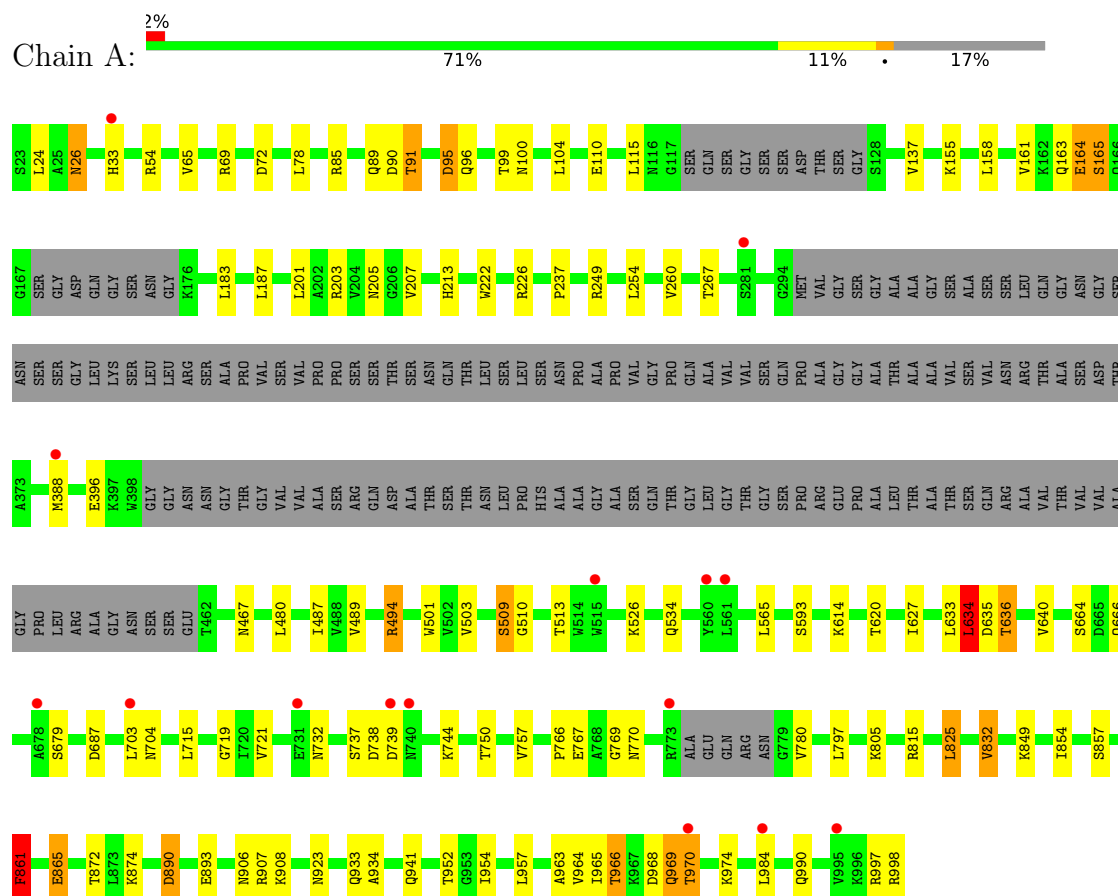
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	54	Total	O	0	0
			54	54		
2	B	10	Total	O	0	0
			10	10		

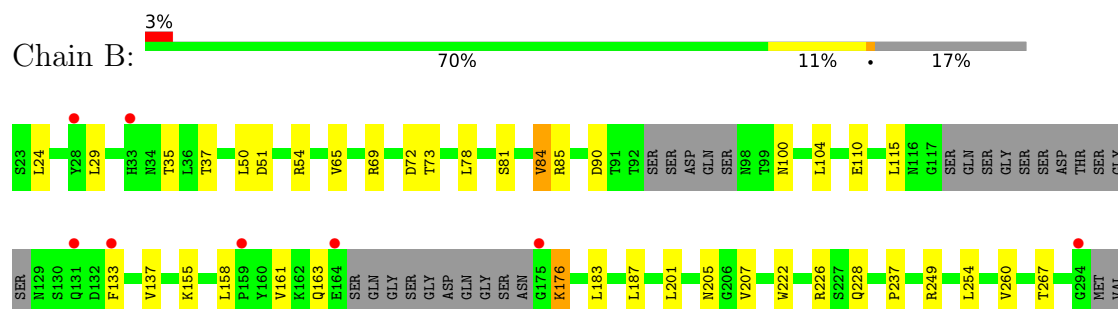
3 Residue-property plots

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: Mgp-operon protein 3



• Molecule 1: Mgp-operon protein 3





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	97.80Å 114.43Å 165.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	82.53 – 2.65 82.53 – 2.65	Depositor EDS
% Data completeness (in resolution range)	99.6 (82.53-2.65) 99.6 (82.53-2.65)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.24 (at 2.65Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.214 , 0.234 0.217 , 0.240	Depositor DCC
R_{free} test set	2655 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	71.7	Xtriage
Anisotropy	0.444	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 59.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12623	wwPDB-VP
Average B, all atoms (Å ²)	89.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.72% 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

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	0.71	0/6422	1.24	18/8731 (0.2%)
1	B	0.70	0/6397	1.21	15/8698 (0.2%)
All	All	0.70	0/12819	1.23	33/17429 (0.2%)

There are no bond length outliers.

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	861	PHE	CA-CB-CG	7.71	121.51	113.80
1	B	861	PHE	CA-CB-CG	7.49	121.29	113.80
1	A	26	ASN	CA-CB-CG	6.97	119.57	112.60
1	B	890	ASP	CA-CB-CG	5.61	118.21	112.60
1	B	907	ARG	N-CA-C	-5.51	106.15	112.92
1	A	90	ASP	CA-C-N	5.49	132.02	121.54
1	A	90	ASP	C-N-CA	5.49	132.02	121.54
1	A	907	ARG	N-CA-C	-5.48	106.17	112.92
1	B	90	ASP	CA-CB-CG	5.48	118.08	112.60
1	B	72	ASP	CA-CB-CG	5.47	118.08	112.60
1	B	679	SER	N-CA-C	5.41	112.88	108.07
1	A	968	ASP	CA-C-N	5.35	131.76	121.54
1	A	968	ASP	C-N-CA	5.35	131.76	121.54
1	A	890	ASP	CA-CB-CG	5.31	117.91	112.60
1	B	702	LYS	CA-C-N	5.27	127.66	120.54
1	B	702	LYS	C-N-CA	5.27	127.66	120.54
1	B	952	THR	N-CA-C	-5.27	106.54	113.12
1	A	72	ASP	CA-CB-CG	5.25	117.85	112.60
1	A	396	GLU	CA-C-N	5.21	128.64	120.82
1	A	396	GLU	C-N-CA	5.21	128.64	120.82
1	A	226	ARG	CA-C-N	5.20	127.56	120.54
1	A	226	ARG	C-N-CA	5.20	127.56	120.54
1	A	679	SER	N-CA-C	5.15	112.66	108.07
1	B	51	ASP	CA-C-N	5.10	127.72	120.53
1	B	51	ASP	C-N-CA	5.10	127.72	120.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	740	ASN	CA-CB-CG	5.10	117.70	112.60
1	B	176	LYS	CA-C-N	5.09	124.80	119.92
1	B	176	LYS	C-N-CA	5.09	124.80	119.92
1	B	493	ASP	CA-CB-CG	5.08	117.68	112.60
1	A	970	THR	CA-C-N	5.07	127.03	120.44
1	A	970	THR	C-N-CA	5.07	127.03	120.44
1	A	164	GLU	CA-C-N	5.02	131.13	121.54
1	A	164	GLU	C-N-CA	5.02	131.13	121.54

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	6292	0	6117	29	0
1	B	6267	0	6091	33	0
2	A	54	0	0	1	0
2	B	10	0	0	0	0
All	All	12623	0	12208	62	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 2.

All (62) 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:775:GLU:C	1:B:776:GLN:HG3	1.81	1.02
1:A:861:PHE:HB2	1:A:865:GLU:HG3	1.56	0.86
1:B:861:PHE:HB2	1:B:865:GLU:HG3	1.56	0.86
1:B:78:LEU:HB2	1:B:480:LEU:HD23	1.65	0.79
1:A:78:LEU:HB2	1:A:480:LEU:HD23	1.64	0.78
1:B:81:SER:O	1:B:84:VAL:HG23	1.85	0.77
1:A:634:LEU:HD22	1:A:635:ASP:H	1.50	0.75
1:A:155:LYS:HB3	1:A:509:SER:HB2	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:155:LYS:HB3	1:B:509:SER:HB2	1.73	0.70
1:B:614:LYS:HD3	1:B:849:LYS:HB3	1.79	0.64
1:B:703:LEU:HD22	1:B:703:LEU:H	1.66	0.61
1:B:537:PHE:HB3	1:B:672:ILE:HD13	1.85	0.58
1:B:775:GLU:C	1:B:776:GLN:CG	2.65	0.56
1:A:766:PRO:HG3	1:A:780:VAL:HB	1.87	0.56
1:A:95:ASP:HA	1:A:165:SER:HB3	1.89	0.54
1:B:775:GLU:O	1:B:776:GLN:HG3	2.07	0.54
1:A:872:THR:HB	1:A:890:ASP:O	2.07	0.53
1:B:719:GLY:HA3	1:B:750:THR:HG22	1.89	0.53
1:A:719:GLY:HA3	1:A:750:THR:HG22	1.90	0.52
1:A:952:THR:HB	1:A:965:ILE:HB	1.90	0.52
1:A:966:THR:HG22	1:A:974:LYS:HB3	1.92	0.52
1:B:205:ASN:HD21	1:B:687:ASP:HB2	1.75	0.52
1:A:636:THR:OG1	1:A:769:GLY:HA2	2.10	0.52
1:A:205:ASN:HD21	1:A:687:ASP:HB2	1.75	0.51
1:A:503:VAL:HG12	1:A:510:GLY:HA3	1.93	0.50
1:A:634:LEU:HD22	1:A:635:ASP:N	2.25	0.50
1:B:503:VAL:HG12	1:B:510:GLY:HA3	1.92	0.50
1:B:825:LEU:HG	1:B:832:VAL:HG13	1.96	0.48
1:A:633:LEU:C	1:A:634:LEU:HD13	2.38	0.48
1:B:954:ILE:HG12	1:B:964:VAL:HG22	1.96	0.48
1:B:957:LEU:HD11	1:B:963:ALA:HB2	1.96	0.48
1:A:825:LEU:HG	1:A:832:VAL:HG13	1.96	0.47
1:A:614:LYS:HD2	1:A:849:LYS:HB2	1.98	0.46
1:B:633:LEU:HG	1:B:634:LEU:HD22	1.97	0.46
1:A:489:VAL:HB	1:A:501:TRP:HB2	1.98	0.46
1:A:957:LEU:HD11	1:A:963:ALA:HB2	1.96	0.46
1:A:513:THR:HG23	1:A:732:ASN:HD22	1.82	0.45
1:B:537:PHE:CB	1:B:672:ILE:HD13	2.45	0.45
1:B:85:ARG:HG3	1:B:110:GLU:HB3	1.98	0.45
1:A:85:ARG:HG3	1:A:110:GLU:HB3	1.99	0.45
1:A:954:ILE:HG12	1:A:964:VAL:HG22	1.99	0.45
1:A:861:PHE:HB2	1:A:865:GLU:CG	2.37	0.44
1:B:778:ASN:HB3	1:B:779:GLY:H	1.61	0.44
1:B:84:VAL:HG22	1:B:133:PHE:O	2.18	0.44
1:B:222:TRP:CD2	1:B:237:PRO:HB3	2.54	0.43
1:B:777:ARG:HE	1:B:777:ARG:HB3	1.50	0.43
1:B:489:VAL:HB	1:B:501:TRP:HB2	1.98	0.43
1:B:775:GLU:O	1:B:776:GLN:CG	2.66	0.43
1:A:222:TRP:CD2	1:A:237:PRO:HB3	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:81:SER:O	1:B:84:VAL:CG2	2.63	0.43
1:B:854:ILE:HA	1:B:934:ALA:HB2	2.01	0.43
1:B:137:VAL:HG11	1:B:487:ILE:HG21	2.00	0.43
1:A:388:MET:HE3	1:A:494:ARG:HG2	2.01	0.42
1:B:861:PHE:HB2	1:B:865:GLU:CG	2.37	0.42
1:A:213:HIS:HE1	2:A:1029:HOH:O	2.02	0.42
1:B:592:ALA:HB3	1:B:595:TYR:HB2	2.02	0.41
1:A:161:VAL:HG21	1:A:183:LEU:HD12	2.03	0.41
1:A:854:ILE:HA	1:A:934:ALA:HB2	2.02	0.41
1:B:991:LEU:O	1:B:995:VAL:HG23	2.21	0.41
1:B:35:THR:HA	1:B:50:LEU:O	2.21	0.40
1:A:137:VAL:HG11	1:A:487:ILE:HG21	2.02	0.40
1:B:161:VAL:HG21	1:B:183:LEU:HD12	2.03	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	802/976 (82%)	751 (94%)	37 (5%)	14 (2%)	7	12
1	B	796/976 (82%)	749 (94%)	38 (5%)	9 (1%)	11	19
All	All	1598/1952 (82%)	1500 (94%)	75 (5%)	23 (1%)	9	14

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	91	THR
1	A	165	SER
1	A	634	LEU
1	A	96	GLN
1	A	164	GLU

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Mol	Chain	Res	Type
1	A	737	SER
1	A	969	GLN
1	A	997	ARG
1	B	736	THR
1	B	737	SER
1	A	467	ASN
1	A	738	ASP
1	A	770	ASN
1	B	467	ASN
1	B	735	SER
1	B	776	GLN
1	A	33[A]	HIS
1	A	33[B]	HIS
1	A	627	ILE
1	B	627	ILE
1	B	634	LEU
1	B	929	THR
1	B	635	ASP

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	684/800 (86%)	620 (91%)	64 (9%)	8	13
1	B	679/800 (85%)	613 (90%)	66 (10%)	8	12
All	All	1363/1600 (85%)	1233 (90%)	130 (10%)	8	13

All (130) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	LEU
1	A	26	ASN
1	A	54[A]	ARG
1	A	54[B]	ARG
1	A	65	VAL

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Mol	Chain	Res	Type
1	A	69	ARG
1	A	89	GLN
1	A	91	THR
1	A	95	ASP
1	A	99	THR
1	A	100	ASN
1	A	104	LEU
1	A	115	LEU
1	A	158	LEU
1	A	163	GLN
1	A	187	LEU
1	A	201	LEU
1	A	203	ARG
1	A	207	VAL
1	A	249	ARG
1	A	254	LEU
1	A	260	VAL
1	A	267	THR
1	A	494	ARG
1	A	509	SER
1	A	526	LYS
1	A	534	GLN
1	A	565	LEU
1	A	593	SER
1	A	620	THR
1	A	634	LEU
1	A	636	THR
1	A	640	VAL
1	A	664	SER
1	A	666	GLN
1	A	703	LEU
1	A	704	ASN
1	A	715	LEU
1	A	721	VAL
1	A	739	ASP
1	A	744	LYS
1	A	757	VAL
1	A	767	GLU
1	A	797	LEU
1	A	805	LYS
1	A	815	ARG
1	A	825	LEU

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Mol	Chain	Res	Type
1	A	832	VAL
1	A	857	SER
1	A	861	PHE
1	A	865	GLU
1	A	874	LYS
1	A	893	GLU
1	A	906	ASN
1	A	908	LYS
1	A	923	ASN
1	A	933	GLN
1	A	941	GLN
1	A	966	THR
1	A	969	GLN
1	A	970	THR
1	A	984	LEU
1	A	990	GLN
1	A	998	ARG
1	B	24	LEU
1	B	29	LEU
1	B	37	THR
1	B	54	ARG
1	B	65	VAL
1	B	69	ARG
1	B	73	THR
1	B	84	VAL
1	B	100	ASN
1	B	104	LEU
1	B	115	LEU
1	B	158	LEU
1	B	163	GLN
1	B	176	LYS
1	B	187	LEU
1	B	201	LEU
1	B	207	VAL
1	B	226	ARG
1	B	228	GLN
1	B	249	ARG
1	B	254	LEU
1	B	260	VAL
1	B	267	THR
1	B	377	LYS
1	B	463	ASP

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Mol	Chain	Res	Type
1	B	509	SER
1	B	526	LYS
1	B	534	GLN
1	B	565	LEU
1	B	568	VAL
1	B	593	SER
1	B	614	LYS
1	B	620	THR
1	B	634	LEU
1	B	640	VAL
1	B	664	SER
1	B	666	GLN
1	B	703	LEU
1	B	715	LEU
1	B	721	VAL
1	B	738	ASP
1	B	739	ASP
1	B	757	VAL
1	B	776	GLN
1	B	777	ARG
1	B	797	LEU
1	B	805	LYS
1	B	815	ARG
1	B	825	LEU
1	B	832	VAL
1	B	850	ASN
1	B	857	SER
1	B	861	PHE
1	B	865	GLU
1	B	893	GLU
1	B	906	ASN
1	B	908	LYS
1	B	933	GLN
1	B	941	GLN
1	B	965	ILE
1	B	969	GLN
1	B	979	LYS
1	B	986	GLU
1	B	990	GLN
1	B	997	ARG
1	B	998	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	166	GLN
1	A	205	ASN
1	A	218	GLN
1	A	387	GLN
1	A	486	GLN
1	A	557	GLN
1	A	732	ASN
1	A	734	GLN
1	A	906	ASN
1	A	923	ASN
1	A	948	GLN
1	B	101	GLN
1	B	205	ASN
1	B	223	ASN
1	B	387	GLN
1	B	486	GLN
1	B	557	GLN
1	B	624	ASN
1	B	732	ASN
1	B	734	GLN
1	B	906	ASN
1	B	969	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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	812/976 (83%)	0.15	15 (1%) 67 63	36, 76, 125, 169	2 (0%)
1	B	806/976 (82%)	0.29	28 (3%) 47 38	36, 93, 140, 177	4 (0%)
All	All	1618/1952 (82%)	0.22	43 (2%) 56 49	36, 83, 135, 177	6 (0%)

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	294	GLY	4.2
1	B	561	LEU	3.8
1	B	493	ASP	3.6
1	B	399	GLY	3.5
1	B	28[A]	TYR	3.3
1	B	159	PRO	3.3
1	B	701	PHE	3.2
1	A	739	ASP	3.2
1	B	708	LEU	3.1
1	B	732	ASN	3.0
1	B	983	GLY	3.0
1	B	633	LEU	2.9
1	B	514	TRP	2.9
1	A	33[A]	HIS	2.7
1	B	972	THR	2.7
1	A	388	MET	2.7
1	A	970	THR	2.7
1	A	773	ARG	2.6
1	A	731	GLU	2.6
1	B	33[A]	HIS	2.6
1	A	678	ALA	2.6
1	A	740	ASN	2.5
1	A	995	VAL	2.5
1	B	930	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	560	TYR	2.4
1	B	774	ALA	2.4
1	A	703	LEU	2.4
1	B	131	GLN	2.3
1	B	396	GLU	2.3
1	B	998	ARG	2.3
1	B	164	GLU	2.3
1	B	758	PRO	2.3
1	B	651	GLN	2.3
1	A	281	SER	2.2
1	B	133	PHE	2.2
1	B	636	THR	2.1
1	A	984	LEU	2.1
1	B	175	GLY	2.1
1	B	512	ALA	2.1
1	A	515	TRP	2.1
1	A	561	LEU	2.1
1	B	847	THR	2.0
1	B	971	TRP	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.