



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 03:47 am BST

PDB ID : 6K68
Title : Application of anti-helix antibodies in protein structure determination (8420-3MNZ)
Authors : Lee, J.O.; Jin, M.S.; Kim, J.W.; Kim, S.; Lee, H.; Cho, G.Y.
Deposited on : 2019-06-01
Resolution : 3.20 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
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

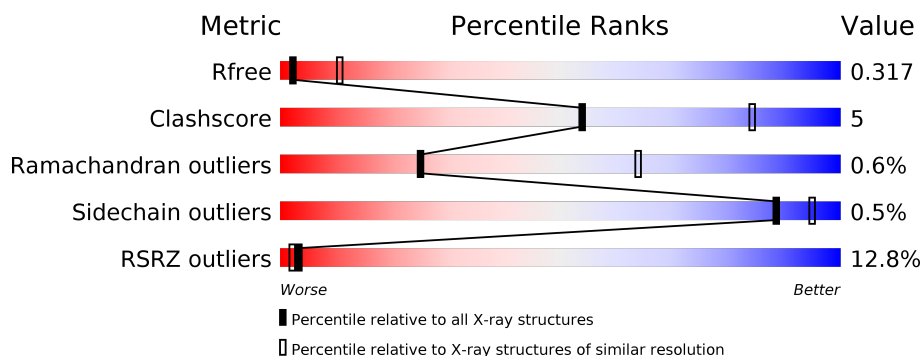
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)

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	115	<div> <div></div> <div>89%</div> <div>7%</div> <div>.</div> </div>
1	C	115	<div> <div></div> <div>88%</div> <div>8%</div> <div>.</div> </div>
1	F	115	<div> <div>5%</div> <div></div> <div>88%</div> <div>8%</div> <div>.</div> </div>
1	J	115	<div> <div>5%</div> <div></div> <div>86%</div> <div>10%</div> <div>..</div> </div>
2	B	116	<div> <div></div> <div>87%</div> <div>9%</div> <div>..</div> </div>
2	D	116	<div> <div>%</div> <div></div> <div>83%</div> <div>13%</div> <div>..</div> </div>

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Mol	Chain	Length	Quality of chain
2	G	116	<div><div></div><div>15%</div><div>84%</div><div>12%</div><div></div><div></div></div>
2	K	116	<div><div></div><div>28%</div><div>83%</div><div>13%</div><div></div><div></div></div>
3	E	73	<div><div></div><div>4%</div><div>59%</div><div>10%</div><div>32%</div><div></div></div>
3	H	73	<div><div></div><div>3%</div><div>62%</div><div>7%</div><div>32%</div><div></div></div>
3	I	73	<div><div></div><div>49%</div><div>56%</div><div>12%</div><div>32%</div><div></div></div>
3	L	73	<div><div></div><div>48%</div><div>33%</div><div>34%</div><div></div><div></div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8562 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 3MNZ Variable heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	110	Total	C	N	O	S	0	0	0
			866	559	138	166	3			
1	C	110	Total	C	N	O	S	0	0	0
			866	559	138	166	3			
1	F	110	Total	C	N	O	S	0	0	0
			866	559	138	166	3			
1	J	111	Total	C	N	O	S	0	0	0
			872	562	139	168	3			

- Molecule 2 is a protein called 3MNZ Variable light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	112	Total	C	N	O	S	0	0	0
			872	548	149	171	4			
2	D	112	Total	C	N	O	S	0	0	0
			872	548	149	171	4			
2	G	112	Total	C	N	O	S	0	0	0
			872	548	149	171	4			
2	K	112	Total	C	N	O	S	0	0	0
			872	548	149	171	4			

- Molecule 3 is a protein called Protein A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	50	Total	C	N	O	S	0	0	0
			401	248	69	83	1			
3	H	50	Total	C	N	O	S	0	0	0
			401	248	69	83	1			
3	I	50	Total	C	N	O	S	0	0	0
			401	248	69	83	1			
3	L	50	Total	C	N	O	S	0	0	0
			401	248	69	83	1			

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.

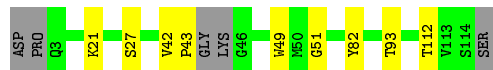
- Molecule 1: 3MNZ Variable heavy chain

Chain A: 




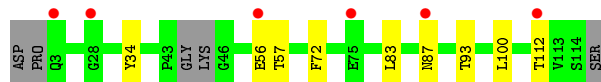
- Molecule 1: 3MNZ Variable heavy chain

Chain C: 




- Molecule 1: 3MNZ Variable heavy chain

Chain F: 




- Molecule 1: 3MNZ Variable heavy chain

Chain J: 

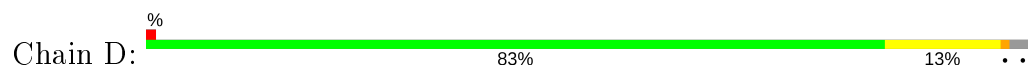


- Molecule 2: 3MNZ Variable light chain

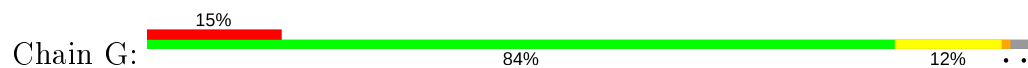
Chain B: 



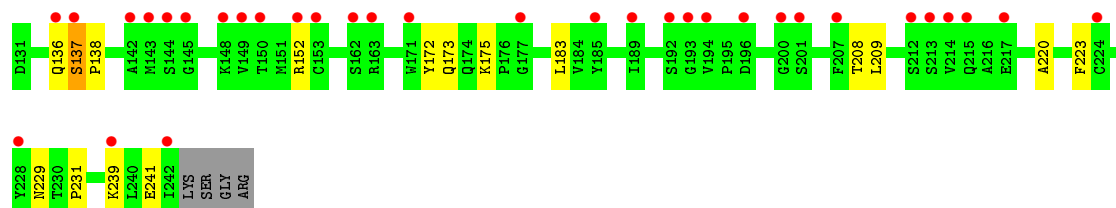
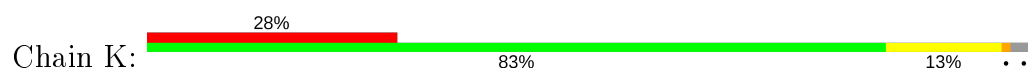
- Molecule 2: 3MNZ Variable light chain



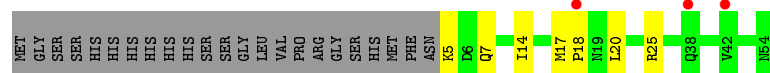
- Molecule 2: 3MNZ Variable light chain



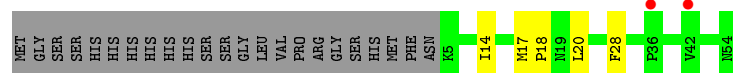
- Molecule 2: 3MNZ Variable light chain



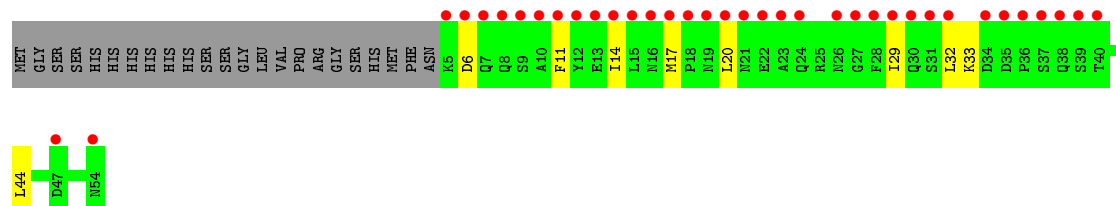
- Molecule 3: Protein A



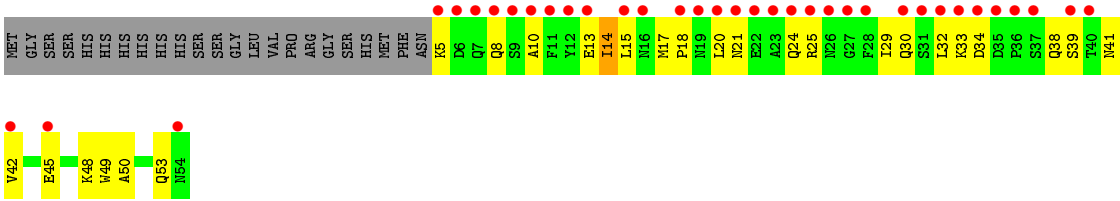
- Molecule 3: Protein A



- Molecule 3: Protein A



- Molecule 3: Protein A



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	75.95Å 95.02Å 179.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.42 – 3.20 33.42 – 3.13	Depositor EDS
% Data completeness (in resolution range)	96.2 (33.42-3.20) 96.2 (33.42-3.13)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.30 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.14_3260, PHENIX 1.14_3260	Depositor
R, R_{free}	0.289 , 0.318 0.287 , 0.317	Depositor DCC
R_{free} test set	2000 reflections (8.79%)	wwPDB-VP
Wilson B-factor (Å ²)	52.8	Xtriage
Anisotropy	0.202	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 28.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.81	EDS
Total number of atoms	8562	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.64% 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](#)

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	0.25	0/891	0.45	0/1213
1	C	0.24	0/891	0.45	0/1213
1	F	0.24	0/891	0.45	0/1213
1	J	0.24	0/897	0.44	0/1221
2	B	0.25	0/891	0.48	0/1205
2	D	0.25	0/891	0.45	0/1205
2	G	0.25	0/891	0.46	0/1205
2	K	0.25	0/891	0.46	0/1205
3	E	0.23	0/407	0.35	0/550
3	H	0.23	0/407	0.34	0/550
3	I	0.23	0/407	0.39	0/550
3	L	0.25	0/407	0.40	0/550
All	All	0.25	0/8762	0.44	0/11880

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 [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	866	0	825	5	0
1	C	866	0	825	6	0
1	F	866	0	825	4	0
1	J	872	0	830	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	872	0	845	7	0
2	D	872	0	845	11	0
2	G	872	0	845	9	0
2	K	872	0	845	10	0
3	E	401	0	381	5	0
3	H	401	0	381	3	0
3	I	401	0	381	7	0
3	L	401	0	381	19	0
All	All	8562	0	8209	89	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 (89) 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)
2:G:137:SER:HB3	2:G:138:PRO:HD3	1.70	0.73
2:K:137:SER:HB3	2:K:138:PRO:HD3	1.72	0.71
2:D:137:SER:HB3	2:D:138:PRO:HD3	1.73	0.70
2:K:136:GLN:HE22	2:K:223:PHE:HA	1.55	0.70
2:G:136:GLN:HE22	2:G:223:PHE:HA	1.61	0.66
2:B:152:ARG:NH2	2:B:206:ASP:OD2	2.30	0.65
1:J:38:TRP:HE1	1:J:53:ILE:HD11	1.60	0.65
3:L:14:ILE:HG21	3:L:29:ILE:HG13	1.76	0.65
2:K:152:ARG:HG2	2:K:208:THR:HG22	1.80	0.63
3:L:38:GLN:HG3	3:L:41:ASN:HB2	1.80	0.63
2:G:167:ASN:HB2	2:G:187:ALA:HB2	1.81	0.61
3:L:14:ILE:HG22	3:L:15:LEU:H	1.66	0.60
2:B:137:SER:HB3	2:B:138:PRO:HD2	1.84	0.58
1:C:93:THR:HG23	1:C:112:THR:HA	1.85	0.58
2:D:173:GLN:HB2	2:D:183:LEU:HD11	1.85	0.58
3:L:32:LEU:HD21	3:L:39:SER:HA	1.87	0.56
2:G:173:GLN:HB2	2:G:183:LEU:HD11	1.88	0.56
3:L:20:LEU:HB2	3:L:25:ARG:HE	1.70	0.56
2:B:137:SER:HB2	2:B:152:ARG:H	1.70	0.56
2:D:152:ARG:HG2	2:D:208:THR:HG22	1.87	0.56
2:B:173:GLN:HB2	2:B:183:LEU:HD11	1.86	0.55
2:K:173:GLN:HB2	2:K:183:LEU:HD11	1.89	0.55
3:L:17:MET:HG3	3:L:20:LEU:HD12	1.87	0.55
1:J:52:TRP:HE1	1:J:61:THR:HB	1.72	0.55
2:D:175:LYS:HG2	2:D:220:ALA:HB2	1.88	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:17:MET:HG2	3:H:20:LEU:HG	1.91	0.53
3:L:32:LEU:HG	3:L:42:VAL:HG21	1.92	0.52
1:A:43:PRO:HD3	1:A:94:ALA:HA	1.92	0.52
1:F:34:TYR:HB2	1:F:100:LEU:HD11	1.93	0.50
3:E:14:ILE:HA	3:E:17:MET:HG2	1.93	0.50
3:H:17:MET:HG3	3:H:18:PRO:HD2	1.92	0.50
2:G:189:ILE:HG21	3:I:44:LEU:HD21	1.94	0.49
1:A:60:PRO:HB3	1:A:72:PHE:HB2	1.93	0.49
1:J:13:LEU:HD23	1:J:112:THR:HB	1.94	0.49
3:L:20:LEU:HD21	3:L:50:ALA:HB2	1.95	0.49
3:L:10:ALA:O	3:L:14:ILE:HG12	2.13	0.49
3:I:11:PHE:CE1	3:I:32:LEU:HB3	2.47	0.49
2:K:137:SER:CB	2:K:138:PRO:HD3	2.43	0.49
3:E:17:MET:HB2	3:E:25:ARG:HH12	1.78	0.49
3:L:15:LEU:O	3:L:25:ARG:NH1	2.46	0.49
3:I:29:ILE:O	3:I:33:LYS:HG2	2.12	0.48
1:J:72:PHE:HE1	1:J:83:LEU:HD13	1.78	0.48
3:L:30:GLN:O	3:L:34:ASP:HB2	2.13	0.48
1:C:42:VAL:HG12	1:C:43:PRO:HD2	1.94	0.48
2:K:172:TYR:HB2	2:K:223:PHE:CE1	2.48	0.48
1:F:56:GLU:HG3	1:F:57:THR:HG23	1.94	0.48
3:L:21:ASN:O	3:L:25:ARG:N	2.36	0.47
2:K:239:LYS:HG2	2:K:241:GLU:OE2	2.14	0.47
3:I:14:ILE:HD12	3:I:32:LEU:HD11	1.96	0.47
2:D:163:ARG:NH1	3:E:18:PRO:HG3	2.29	0.47
3:L:21:ASN:HB3	3:L:24:GLN:HB3	1.96	0.46
3:L:8:GLN:O	3:L:10:ALA:N	2.44	0.46
3:L:10:ALA:O	3:L:13:GLU:N	2.47	0.46
2:G:137:SER:CB	2:G:138:PRO:HD3	2.43	0.46
2:B:167:ASN:HB2	2:B:187:ALA:HB2	1.98	0.46
2:B:175:LYS:HG2	2:B:220:ALA:HB2	1.98	0.45
2:D:137:SER:CB	2:D:138:PRO:HD3	2.42	0.45
3:E:17:MET:HG3	3:E:20:LEU:HD12	1.99	0.45
2:G:175:LYS:HG2	2:G:220:ALA:HB2	1.99	0.45
3:L:49:TRP:O	3:L:53:GLN:HG3	2.17	0.45
2:D:239:LYS:HG2	2:D:241:GLU:OE2	2.17	0.45
3:I:11:PHE:HE1	3:I:32:LEU:HB3	1.82	0.44
1:F:72:PHE:HE1	1:F:83:LEU:HD13	1.82	0.44
3:I:17:MET:HG2	3:I:20:LEU:HG	1.99	0.44
1:A:21:LYS:HE2	1:A:82:TYR:CD1	2.53	0.44
1:F:93:THR:HG23	1:F:112:THR:HA	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:137:SER:HB2	2:D:152:ARG:HB2	2.00	0.43
2:K:209:LEU:HA	2:K:209:LEU:HD23	1.75	0.43
2:G:137:SER:HB2	2:G:152:ARG:HB2	1.99	0.43
3:E:5:LYS:O	3:E:7:GLN:N	2.42	0.42
1:C:21:LYS:HE2	1:C:82:TYR:CD1	2.54	0.42
2:K:175:LYS:HG2	2:K:220:ALA:HB2	2.00	0.42
2:G:227:HIS:HB2	2:G:232:TRP:CZ2	2.54	0.42
1:A:30:THR:OG1	1:C:27:SER:HA	2.19	0.42
1:J:39:VAL:HG21	1:J:105:TRP:CH2	2.55	0.41
1:C:49:TRP:CG	2:D:232:TRP:HB2	2.56	0.41
3:I:20:LEU:HA	3:I:20:LEU:HD23	1.95	0.41
1:J:21:LYS:HE2	1:J:82:TYR:CD1	2.56	0.41
3:L:29:ILE:O	3:L:33:LYS:HB3	2.21	0.41
2:D:148:LYS:HG3	2:D:212:SER:HA	2.03	0.41
3:L:45:GLU:HA	3:L:48:LYS:HB3	2.03	0.41
2:B:152:ARG:HH12	2:B:154:LYS:HE3	1.85	0.41
1:J:52:TRP:NE1	1:J:61:THR:HB	2.34	0.41
1:A:72:PHE:HE1	1:A:83:LEU:HD13	1.85	0.40
3:L:5:LYS:HA	3:L:8:GLN:OE1	2.21	0.40
2:K:229:ASN:HB3	2:K:231:PRO:HD2	2.03	0.40
1:C:49:TRP:CH2	1:C:51:GLY:HA2	2.56	0.40
2:D:225:LEU:HD13	2:D:234:PHE:CE1	2.56	0.40
3:H:14:ILE:HD13	3:H:28:PHE:HB3	2.04	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	106/115 (92%)	100 (94%)	6 (6%)	0	100	100
1	C	106/115 (92%)	100 (94%)	6 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	106/115 (92%)	101 (95%)	5 (5%)	0	100	100
1	J	107/115 (93%)	102 (95%)	5 (5%)	0	100	100
2	B	110/116 (95%)	104 (94%)	5 (4%)	1 (1%)	17	56
2	D	110/116 (95%)	105 (96%)	4 (4%)	1 (1%)	17	56
2	G	110/116 (95%)	105 (96%)	4 (4%)	1 (1%)	17	56
2	K	110/116 (95%)	102 (93%)	7 (6%)	1 (1%)	17	56
3	E	48/73 (66%)	46 (96%)	2 (4%)	0	100	100
3	H	48/73 (66%)	47 (98%)	1 (2%)	0	100	100
3	I	48/73 (66%)	45 (94%)	3 (6%)	0	100	100
3	L	48/73 (66%)	39 (81%)	7 (15%)	2 (4%)	3	20
All	All	1057/1216 (87%)	996 (94%)	55 (5%)	6 (1%)	25	64

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	137	SER
2	D	137	SER
2	G	137	SER
2	K	137	SER
3	L	18	PRO
3	L	14	ILE

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	93/97 (96%)	93 (100%)	0	100	100
1	C	93/97 (96%)	93 (100%)	0	100	100
1	F	93/97 (96%)	92 (99%)	1 (1%)	73	88
1	J	94/97 (97%)	93 (99%)	1 (1%)	73	88
2	B	97/100 (97%)	96 (99%)	1 (1%)	76	90

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	D	97/100 (97%)	97 (100%)	0	100	100
2	G	97/100 (97%)	96 (99%)	1 (1%)	76	90
2	K	97/100 (97%)	97 (100%)	0	100	100
3	E	45/65 (69%)	45 (100%)	0	100	100
3	H	45/65 (69%)	45 (100%)	0	100	100
3	I	45/65 (69%)	44 (98%)	1 (2%)	52	79
3	L	45/65 (69%)	45 (100%)	0	100	100
All	All	941/1048 (90%)	936 (100%)	5 (0%)	88	95

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	137	SER
1	F	87	ASN
2	G	226	GLN
3	I	6	ASP
1	J	52	TRP

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

Mol	Chain	Res	Type
3	E	7	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

There are no ligands in this entry.

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, 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	110/115 (95%)	-0.05	0	100	100	14, 24, 43, 47	0
1	C	110/115 (95%)	0.08	0	100	100	18, 27, 50, 55	0
1	F	110/115 (95%)	0.69	6 (5%)	25	14	49, 73, 86, 94	0
1	J	111/115 (96%)	0.70	6 (5%)	25	14	45, 76, 95, 100	0
2	B	112/116 (96%)	0.16	0	100	100	22, 33, 47, 54	0
2	D	112/116 (96%)	0.19	1 (0%)	84	75	25, 39, 50, 59	1 (0%)
2	G	112/116 (96%)	1.02	17 (15%)	2	1	79, 90, 117, 120	0
2	K	112/116 (96%)	1.65	33 (29%)	0	0	85, 105, 120, 126	0
3	E	50/73 (68%)	0.35	3 (6%)	21	12	23, 42, 72, 75	0
3	H	50/73 (68%)	0.46	2 (4%)	38	25	21, 40, 64, 74	0
3	I	50/73 (68%)	3.13	36 (72%)	0	0	90, 119, 130, 134	0
3	L	50/73 (68%)	3.80	35 (70%)	0	0	98, 129, 136, 138	0
All	All	1089/1216 (89%)	0.81	139 (12%)	3	2	14, 60, 122, 138	1 (0%)

All (139) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	L	23	ALA	11.8
3	L	24	GLN	11.2
3	L	28	PHE	8.8
3	L	19	ASN	8.3
3	L	21	ASN	8.0
3	L	6	ASP	7.6
3	L	27	GLY	7.6
3	L	20	LEU	7.4
3	I	5	LYS	7.0
3	I	31	SER	6.8
3	L	34	ASP	6.8

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Mol	Chain	Res	Type	RSRZ
3	L	9	SER	6.4
3	I	6	ASP	6.2
2	K	162	SER	6.0
3	L	8	GLN	5.5
3	I	19	ASN	5.3
3	L	26	ASN	5.3
3	I	23	ALA	5.2
3	I	18	PRO	5.1
3	I	38	GLN	5.1
2	K	142	ALA	5.1
3	I	28	PHE	5.1
3	I	30	GLN	4.9
3	I	34	ASP	4.9
3	I	12	TYR	4.9
3	I	13	GLU	4.9
3	L	36	PRO	4.9
3	I	35	ASP	4.9
3	I	8	GLN	4.8
2	K	143	MET	4.6
3	I	27	GLY	4.6
3	L	5	LYS	4.5
3	L	33	LYS	4.5
3	I	16	ASN	4.5
2	G	241	GLU	4.4
3	I	37	SER	4.3
3	L	15	LEU	4.3
3	L	12	TYR	4.3
3	I	9	SER	4.1
2	K	153	CYS	4.1
3	L	13	GLU	4.1
3	I	20	LEU	4.1
3	L	16	ASN	4.0
3	L	31	SER	4.0
3	L	30	GLN	3.8
2	G	153	CYS	3.8
2	K	212	SER	3.8
2	G	242	ILE	3.7
2	K	201	SER	3.7
1	J	13	LEU	3.7
3	I	29	ILE	3.6
3	I	17	MET	3.6
3	L	37	SER	3.5

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Mol	Chain	Res	Type	RSRZ
2	K	144	SER	3.5
2	K	196	ASP	3.5
3	I	22	GLU	3.5
3	L	32	LEU	3.4
2	K	228	TYR	3.4
3	I	14	ILE	3.3
2	K	152	ARG	3.3
3	L	42	VAL	3.3
3	I	32	LEU	3.2
2	G	139	SER	3.2
3	I	36	PRO	3.2
3	L	25	ARG	3.2
2	K	150	THR	3.2
3	I	21	ASN	3.2
2	G	146	GLY	3.1
2	G	143	MET	3.1
3	L	7	GLN	3.1
2	K	137	SER	3.0
3	E	18	PRO	3.0
2	G	131	ASP	3.0
2	K	213	SER	3.0
1	F	112	THR	3.0
2	K	192	SER	3.0
3	L	39	SER	3.0
1	J	56	GLU	3.0
1	J	112	THR	2.9
1	J	90	ASN	2.9
2	K	242	ILE	2.9
3	I	54	ASN	2.9
2	G	240	LEU	2.9
2	K	214	VAL	2.9
3	I	11	PHE	2.9
3	L	10	ALA	2.8
2	G	150	THR	2.8
3	L	18	PRO	2.8
2	K	145	GLY	2.8
2	K	194	VAL	2.8
2	K	149	VAL	2.7
2	G	142	ALA	2.7
2	G	205	THR	2.6
1	F	75	GLU	2.6
3	I	26	ASN	2.6

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Mol	Chain	Res	Type	RSRZ
2	K	177	GLY	2.6
2	K	189	ILE	2.6
2	G	140	SER	2.6
2	K	185	TYR	2.5
3	I	15	LEU	2.5
1	F	3	GLN	2.5
2	K	148	LYS	2.5
2	K	136	GLN	2.5
2	D	241	GLU	2.5
2	K	171	TRP	2.5
1	J	43	PRO	2.5
2	K	193	GLY	2.4
2	K	217	GLU	2.4
1	J	64	ASP	2.4
3	I	47	ASP	2.4
1	F	87	ASN	2.4
3	I	40	THR	2.3
1	F	28	GLY	2.3
3	L	11	PHE	2.3
3	I	7	GLN	2.3
2	K	200	GLY	2.3
3	L	45	GLU	2.3
3	I	39	SER	2.3
3	L	40	THR	2.3
3	H	36	PRO	2.3
3	L	22	GLU	2.2
2	G	206	ASP	2.2
2	G	148	LYS	2.2
2	K	239	LYS	2.2
3	L	54	ASN	2.2
2	K	163	ARG	2.2
2	G	192	SER	2.2
2	G	223	PHE	2.1
2	K	224	CYS	2.1
3	H	42	VAL	2.1
3	L	35	ASP	2.1
2	K	215	GLN	2.1
3	I	10	ALA	2.1
3	I	24	GLN	2.1
3	E	42	VAL	2.1
1	F	56	GLU	2.0
2	K	207	PHE	2.0

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Mol	Chain	Res	Type	RSRZ
2	G	136	GLN	2.0
3	E	38	GLN	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.