



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 04:47 PM UTC

PDB ID : 3QJH / pdb_00003qjh
Title : The crystal structure of the 5c.c7 TCR
Authors : Ely, L.K.; Newell, E.W.; Davis, M.M.; Garcia, K.C.
Deposited on : 2011-01-28
Resolution : 1.90 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

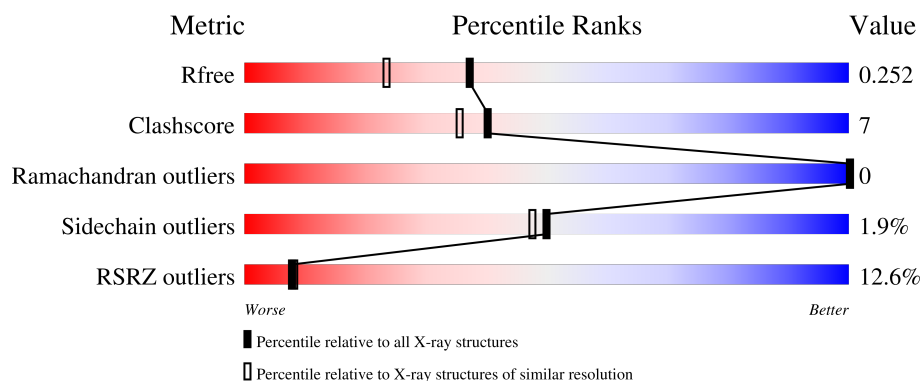
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	205	13% 76% 18% • 5%
1	C	205	15% 83% 12% 5%
2	B	243	12% 87% 12% •
2	D	243	9% 86% 13% •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7223 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 5c.c7 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	195	Total	C	N	O	S	0	0	0
			1502	933	259	302	8			
1	C	195	Total	C	N	O	S	0	0	0
			1503	935	259	301	8			

- Molecule 2 is a protein called 5c.c7 beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	243	Total	C	N	O	S	0	0	0
			1915	1206	333	367	9			
2	D	242	Total	C	N	O	S	0	0	0
			1915	1208	334	364	9			

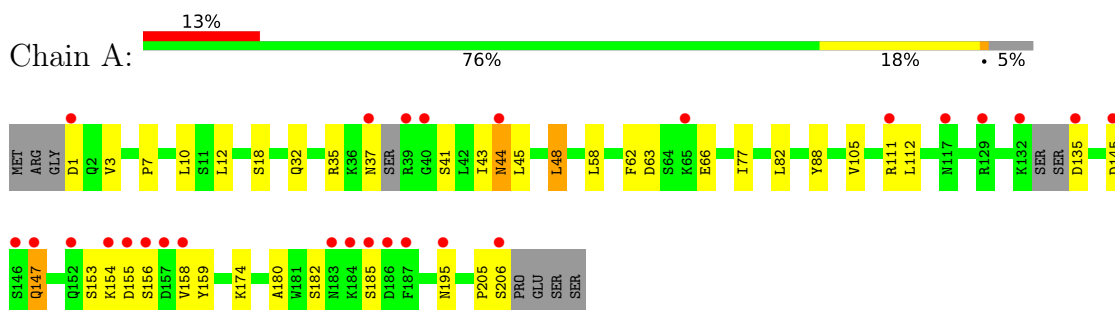
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	78	Total	O	0	0
			78	78		
3	B	111	Total	O	0	0
			111	111		
3	C	79	Total	O	0	0
			79	79		
3	D	120	Total	O	0	0
			120	120		

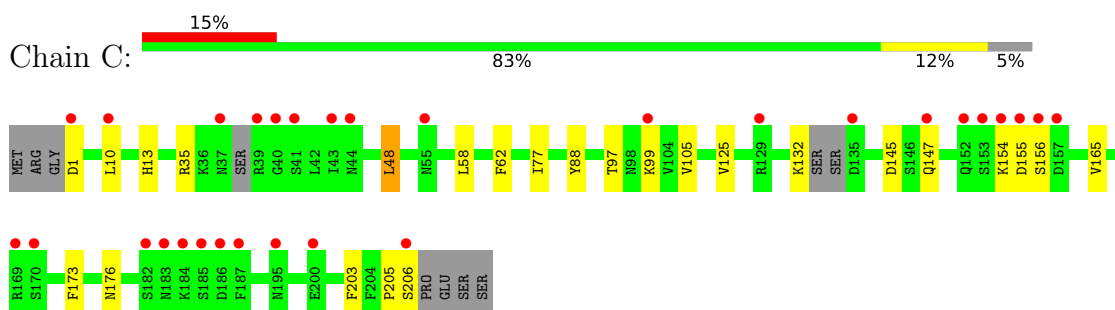
3 Residue-property plots [i](#)

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.

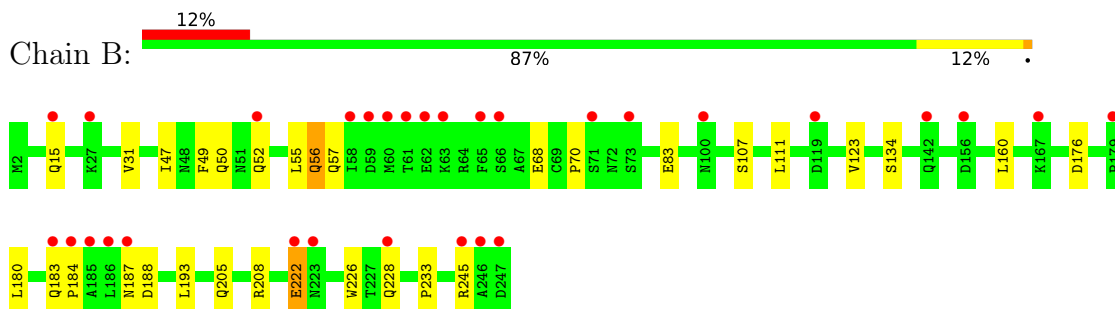
- Molecule 1: 5c.c7 alpha chain



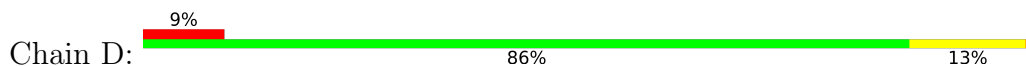
- Molecule 1: 5c.c7 alpha chain

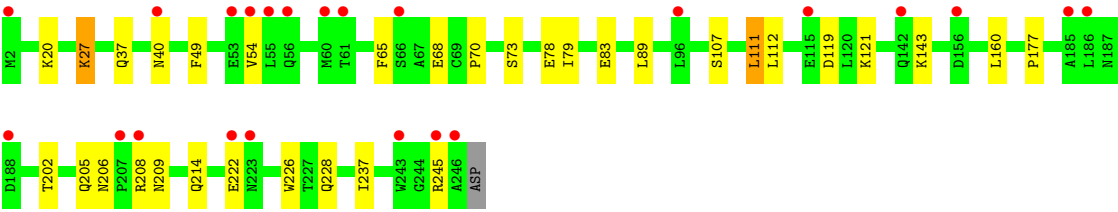


- Molecule 2: 5c.c7 beta chain



- Molecule 2: 5c.c7 beta chain





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	133.17Å 139.46Å 61.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.32 – 1.90 46.32 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.5 (46.32-1.90) 99.5 (46.32-1.90)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.35 (at 1.90Å)	Xtriage
Refinement program	PHENIX 1.6.4_486	Depositor
R, R_{free}	0.225 , 0.256 0.223 , 0.252	Depositor DCC
R_{free} test set	4568 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	28.5	Xtriage
Anisotropy	0.237	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 38.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.011 for k,h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7223	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.27% 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	0.45	0/1529	0.76	0/2069
1	C	0.48	0/1530	0.74	0/2069
2	B	0.47	0/1964	0.74	2/2673 (0.1%)
2	D	0.49	0/1964	0.74	0/2672
All	All	0.47	0/6987	0.74	2/9483 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	183	GLN	CA-C-N	5.84	125.32	119.24
2	B	183	GLN	C-N-CA	5.84	125.32	119.24

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	1502	0	1413	31	0
1	C	1503	0	1420	20	0
2	B	1915	0	1795	20	1
2	D	1915	0	1816	26	1
3	A	78	0	0	3	0
3	B	111	0	0	4	0
3	C	79	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	120	0	0	4	0
All	All	7223	0	6444	88	2

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 7.

All (88) 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:154:LYS:HD3	1:C:156:SER:HA	1.50	0.94
1:A:32:GLN:NE2	1:A:44:ASN:OD1	2.04	0.90
2:B:184:PRO:HG2	1:C:154:LYS:HG3	1.59	0.84
2:B:15:GLN:HE22	2:B:83:GLU:HG2	1.44	0.81
1:A:1:ASP:O	1:A:105:VAL:HG23	1.86	0.75
1:C:97:THR:OG1	1:C:99:LYS:HG2	1.88	0.74
2:B:52:GLN:OE1	3:B:331:HOH:O	2.08	0.71
2:B:15:GLN:NE2	2:B:83:GLU:HG2	2.05	0.71
3:A:281:HOH:O	1:C:13:HIS:HE1	1.76	0.67
1:A:154:LYS:CD	1:C:156:SER:HA	2.25	0.66
2:D:222:GLU:HA	3:D:295:HOH:O	1.97	0.65
1:C:132:LYS:HE2	3:C:245:HOH:O	1.97	0.65
1:A:185:SER:HA	3:A:422:HOH:O	1.98	0.64
1:A:147:GLN:OE1	2:D:40:ASN:CG	2.41	0.62
1:A:35:ARG:HD3	1:A:88:TYR:CZ	2.34	0.62
2:D:65:PHE:CE2	2:D:79:ILE:HG12	2.34	0.62
1:A:153:SER:CB	1:A:158:VAL:HG23	2.30	0.61
2:B:56:GLN:HG2	3:B:315:HOH:O	2.01	0.61
1:A:154:LYS:HD3	1:C:156:SER:CA	2.27	0.61
1:A:153:SER:OG	1:A:158:VAL:HG23	2.02	0.60
2:D:222:GLU:N	2:D:222:GLU:OE1	2.35	0.60
2:D:119:ASP:OD1	2:D:121:LYS:HD3	2.03	0.59
2:B:50:GLN:HB2	2:B:55:LEU:HD11	1.84	0.58
2:D:49:PHE:CE1	2:D:54:VAL:HG22	2.40	0.56
1:C:145:ASP:OD2	1:C:147:GLN:HG2	2.05	0.56
1:C:205:PRO:O	1:C:206:SER:HB3	2.05	0.55
2:B:47:ILE:HG13	2:B:57:GLN:HB3	1.88	0.55
2:B:222:GLU:HA	3:B:282:HOH:O	2.06	0.54
1:A:153:SER:HA	1:A:195:ASN:ND2	2.23	0.54
1:C:203:PHE:CE2	1:C:205:PRO:HG3	2.44	0.53
1:A:205:PRO:O	1:A:206:SER:HB3	2.09	0.52
2:D:107:SER:HB2	3:D:270:HOH:O	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:ASP:OD1	1:A:158:VAL:HG22	2.11	0.51
2:B:226:TRP:CE2	2:B:228:GLN:HB2	2.46	0.51
2:B:176:ASP:HB2	2:B:193:LEU:HD12	1.93	0.50
2:D:214:GLN:HG3	2:D:237:ILE:CG2	2.42	0.50
2:D:226:TRP:CZ2	2:D:228:GLN:HG3	2.47	0.50
1:A:82:LEU:HD22	1:A:174:LYS:HD3	1.95	0.49
2:B:180:LEU:C	2:B:180:LEU:HD12	2.38	0.49
1:A:3:VAL:HG23	1:A:105:VAL:HG22	1.94	0.49
2:B:226:TRP:CZ2	2:B:228:GLN:HG3	2.48	0.48
2:D:20:LYS:HE3	2:D:78:GLU:OE2	2.14	0.48
1:C:35:ARG:HD3	1:C:88:TYR:CZ	2.48	0.48
1:A:58:LEU:CD2	1:A:77:ILE:HG12	2.44	0.47
1:C:58:LEU:HD22	1:C:77:ILE:HG12	1.96	0.47
2:D:205:GLN:HA	2:D:245:ARG:O	2.13	0.47
2:D:206:ASN:HB3	2:D:209:ASN:ND2	2.28	0.47
1:A:154:LYS:O	1:C:155:ASP:HA	2.15	0.47
2:B:31:VAL:HA	2:B:49:PHE:O	2.15	0.46
2:D:226:TRP:NE1	2:D:228:GLN:HB2	2.31	0.46
1:A:44:ASN:HD22	1:A:45:LEU:N	2.13	0.46
1:A:48:LEU:HD13	1:A:62:PHE:HB3	1.98	0.46
1:A:159:TYR:O	1:A:180:ALA:HA	2.16	0.46
2:D:111:LEU:HD22	2:D:112:LEU:N	2.31	0.46
2:B:68:GLU:HG3	2:B:70:PRO:HD3	1.98	0.46
2:B:226:TRP:NE1	2:B:228:GLN:HB2	2.31	0.46
1:C:173:PHE:CD2	2:D:143:LYS:HE2	2.51	0.45
1:A:145:ASP:C	1:A:147:GLN:H	2.24	0.45
1:C:165:VAL:HG22	1:C:176:ASN:OD1	2.16	0.45
2:D:37:GLN:HE21	2:D:89:LEU:HD23	1.82	0.45
1:A:145:ASP:CG	1:A:147:GLN:HB2	2.41	0.45
1:A:145:ASP:OD1	1:A:147:GLN:HB2	2.17	0.44
2:D:226:TRP:CE2	2:D:228:GLN:HB2	2.52	0.44
2:B:123:VAL:HG12	2:B:233:PRO:HB2	1.99	0.43
1:C:58:LEU:CD2	1:C:77:ILE:HG12	2.48	0.43
2:B:205:GLN:HA	2:B:245:ARG:O	2.19	0.43
1:A:12:LEU:HD21	1:A:18:SER:HB2	2.01	0.43
1:A:135:ASP:HA	3:A:216:HOH:O	2.19	0.43
1:C:125:VAL:O	1:C:205:PRO:HD2	2.19	0.43
1:A:7:PRO:HB3	1:C:10:LEU:HD11	2.01	0.42
1:A:147:GLN:HE22	2:D:40:ASN:HB3	1.85	0.42
2:D:202:THR:HG23	3:D:343:HOH:O	2.20	0.42
2:B:107:SER:HB2	3:B:272:HOH:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:187:ASN:O	2:B:188:ASP:HB2	2.21	0.41
2:D:160:LEU:C	2:D:160:LEU:HD23	2.45	0.41
2:B:160:LEU:C	2:B:160:LEU:HD23	2.46	0.41
1:A:158:VAL:HG12	1:A:182:SER:HB2	2.03	0.41
2:D:37:GLN:NE2	2:D:89:LEU:HD23	2.36	0.41
1:A:10:LEU:HD23	1:A:112:LEU:HD13	2.02	0.41
1:A:111:ARG:NH1	1:A:112:LEU:O	2.54	0.41
2:D:70:PRO:HB2	2:D:73:SER:HB2	2.03	0.41
2:D:177:PRO:HD2	3:D:360:HOH:O	2.20	0.41
2:D:27:LYS:HB3	2:D:27:LYS:HE2	1.69	0.40
2:D:68:GLU:HG3	2:D:70:PRO:HD3	2.03	0.40
1:C:1:ASP:O	1:C:105:VAL:HG23	2.21	0.40
1:C:48:LEU:HD13	1:C:62:PHE:HB3	2.02	0.40
2:D:214:GLN:HG3	2:D:237:ILE:HG21	2.03	0.40
1:A:63:ASP:CG	1:A:66:GLU:HB2	2.46	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:83:GLU:OE2	2:D:208:ARG:NH1[1_556]	1.90	0.30
2:B:83:GLU:OE2	2:B:208:ARG:NH1[1_554]	2.03	0.17

5.3 Torsion angles

5.3.1 Protein backbone

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	189/205 (92%)	189 (100%)	0	0	100	100
1	C	189/205 (92%)	187 (99%)	2 (1%)	0	100	100
2	B	241/243 (99%)	235 (98%)	6 (2%)	0	100	100
2	D	240/243 (99%)	232 (97%)	8 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	859/896 (96%)	843 (98%)	16 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	166/181 (92%)	159 (96%)	7 (4%)	26	19
1	C	166/181 (92%)	165 (99%)	1 (1%)	78	81
2	B	206/215 (96%)	202 (98%)	4 (2%)	50	47
2	D	208/215 (97%)	206 (99%)	2 (1%)	68	70
All	All	746/792 (94%)	732 (98%)	14 (2%)	50	47

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	ASN
1	A	41	SER
1	A	43	ILE
1	A	44	ASN
1	A	48	LEU
1	A	147	GLN
1	A	156	SER
2	B	56	GLN
2	B	111	LEU
2	B	134	SER
2	B	222	GLU
1	C	48	LEU
2	D	27	LYS
2	D	111	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	127	GLN
1	A	195	ASN
2	B	15	GLN
2	B	36	GLN
2	B	80	GLN
2	B	142	GLN
1	C	147	GLN
2	D	37	GLN
2	D	51	ASN
2	D	80	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](#)

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	195/205 (95%)	0.74	27 (13%) 6 6	21, 31, 49, 67	0
1	C	195/205 (95%)	0.87	30 (15%) 5 5	20, 31, 51, 70	0
2	B	243/243 (100%)	0.71	30 (12%) 8 9	18, 30, 49, 68	2 (0%)
2	D	242/243 (99%)	0.61	23 (9%) 14 14	15, 29, 45, 56	2 (0%)
All	All	875/896 (97%)	0.72	110 (12%) 8 8	15, 30, 49, 70	4 (0%)

All (110) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	61	THR	6.1
1	A	206	SER	5.7
1	C	185	SER	5.0
2	B	186	LEU	5.0
2	B	185	ALA	4.9
2	B	246	ALA	4.9
1	A	185	SER	4.9
1	C	206	SER	4.9
1	C	186	ASP	4.9
2	D	246	ALA	4.7
1	C	37	ASN	4.6
2	D	185	ALA	4.2
1	A	156	SER	4.2
1	C	154	LYS	4.0
1	C	1	ASP	3.9
2	D	186	LEU	3.9
1	A	155	ASP	3.8
2	D	223	ASN	3.8
1	A	117	ASN	3.7
1	C	195	ASN	3.7
1	C	135	ASP	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	184	LYS	3.7
1	C	184	LYS	3.6
1	C	156	SER	3.5
1	A	40	GLY	3.5
2	D	222	GLU	3.5
1	C	147	GLN	3.4
1	C	39	ARG	3.4
1	A	132	LYS	3.4
2	B	156	ASP	3.4
2	D	2	MET	3.3
1	A	39	ARG	3.3
1	A	147	GLN	3.3
2	B	187	ASN	3.3
1	A	1	ASP	3.2
1	C	155	ASP	3.2
1	A	183	ASN	3.2
1	C	157	ASP	3.2
1	C	183	ASN	3.2
2	D	142	GLN	3.2
1	A	186	ASP	3.1
2	D	53	GLU	3.1
1	A	146	SER	3.1
2	B	247	ASP	3.1
2	B	15	GLN	3.1
1	C	200	GLU	3.0
2	B	73	SER	3.0
2	D	96	LEU	3.0
2	D	208	ARG	3.0
2	B	184	PRO	3.0
1	A	37	ASN	3.0
2	B	60	MET	2.9
1	C	129	ARG	2.9
2	B	245	ARG	2.9
2	B	223	ASN	2.9
1	A	135	ASP	2.9
2	B	58	ILE	2.8
1	A	157	ASP	2.8
2	B	222	GLU	2.8
1	A	152	GLN	2.7
1	C	152	GLN	2.7
2	B	142	GLN	2.7
1	C	43	ILE	2.7

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Mol	Chain	Res	Type	RSRZ
2	B	228	GLN	2.7
2	B	63	LYS	2.7
1	A	195	ASN	2.6
2	D	54	VAL	2.6
2	D	115	GLU	2.6
2	D	40	ASN	2.6
1	C	40	GLY	2.6
2	D	61	THR	2.6
1	C	10	LEU	2.6
2	D	56	GLN	2.6
2	B	59	ASP	2.6
1	C	99	LYS	2.6
2	B	65	PHE	2.5
2	B	71	SER	2.5
2	D	188	ASP	2.5
1	C	187	PHE	2.5
1	A	158	VAL	2.5
2	B	167	LYS	2.5
2	B	62	GLU	2.4
1	A	145	ASP	2.4
1	A	111	ARG	2.4
1	A	187	PHE	2.4
2	B	179	PRO	2.4
2	D	207	PRO	2.4
1	C	153	SER	2.4
2	B	66	SER	2.3
1	C	169	ARG	2.3
2	D	66	SER	2.3
2	B	119	ASP	2.3
2	D	60	MET	2.3
2	B	27	LYS	2.3
1	C	55	ASN	2.2
1	C	170	SER	2.2
2	B	52	GLN	2.2
2	D	156	ASP	2.2
1	A	129	ARG	2.2
2	D	55	LEU	2.1
2	D	245	ARG	2.1
1	A	44	ASN	2.1
1	C	44	ASN	2.1
2	B	183	GLN	2.1
1	C	182	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	65	LYS	2.0
1	C	41	SER	2.0
2	D	243	TRP	2.0
1	A	154	LYS	2.0
2	B	100	ASN	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.