



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 11:10 PM EDT

PDB ID : 6DD9 / pdb_00006dd9
Title : Structure of mouse SYCP3, P1 form
Authors : Rosenberg, S.C.; Munoz, I.C.; Uson, I.; Corbett, K.D.
Deposited on : 2018-05-09
Resolution : 2.30 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 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

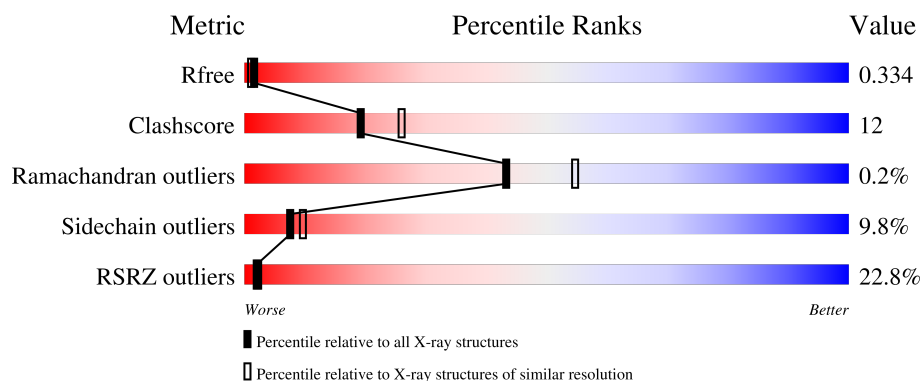
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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 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	144	18% 58% 29% 11%
1	B	144	17% 58% 27% 12%
1	C	144	17% 62% 22% 13%
1	D	144	22% 58% 24% 16%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4119 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 Synaptonemal complex protein 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	128	Total	C	N	O	Se	0	0	0
			1059	662	185	206	6			
1	B	127	Total	C	N	O	Se	0	0	0
			1052	660	184	203	5			
1	D	121	Total	C	N	O	Se	0	0	0
			972	607	173	188	4			
1	C	125	Total	C	N	O	Se	0	0	0
			1026	649	179	192	6			

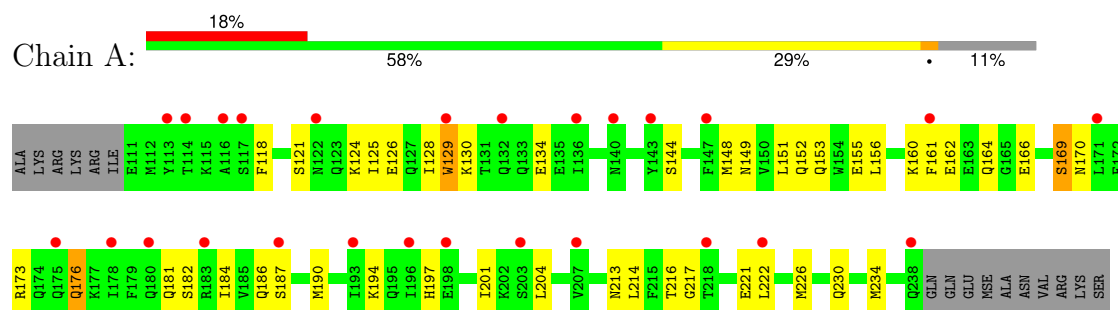
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	4	Total	O	0	0
			4	4		
2	D	4	Total	O	0	0
			4	4		
2	C	2	Total	O	0	0
			2	2		

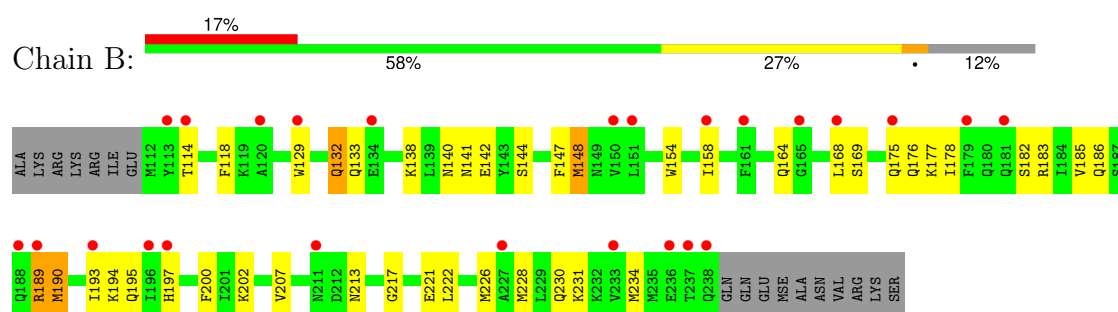
3 Residue-property plots [i](#)

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.

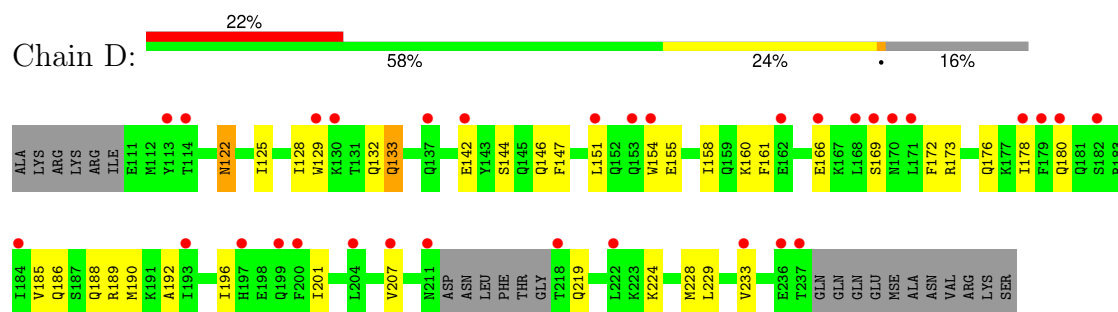
• Molecule 1: Synaptonemal complex protein 3



• Molecule 1: Synaptonemal complex protein 3

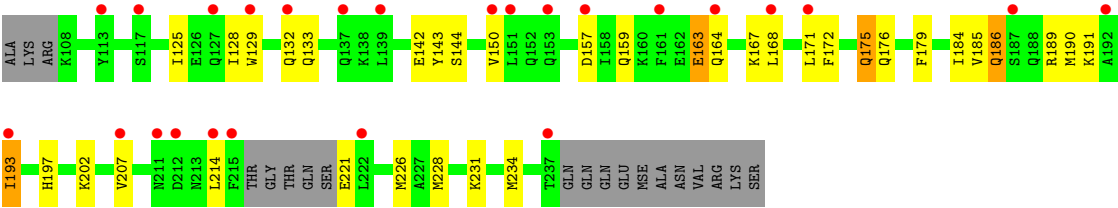


• Molecule 1: Synaptonemal complex protein 3



• Molecule 1: Synaptonemal complex protein 3





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	45.87Å 52.27Å 75.59Å 95.17° 103.70° 110.54°	Depositor
Resolution (Å)	72.10 – 2.30 72.10 – 2.30	Depositor EDS
% Data completeness (in resolution range)	91.5 (72.10-2.30) 94.0 (72.10-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.82 (at 2.29Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.287 , 0.331 0.287 , 0.334	Depositor DCC
R_{free} test set	1307 reflections (4.69%)	wwPDB-VP
Wilson B-factor (Å ²)	50.2	Xtriage
Anisotropy	0.532	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 52.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4119	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.64% 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	0.39	0/1064	0.61	0/1408
1	B	0.37	0/1059	0.56	0/1404
1	C	0.40	0/1030	0.57	0/1360
1	D	0.40	0/975	0.62	0/1292
All	All	0.39	0/4128	0.59	0/5464

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	1059	0	1040	35	0
1	B	1052	0	1029	45	0
1	C	1026	0	1010	33	0
1	D	972	0	919	40	0
2	A	4	0	0	0	0
2	C	2	0	0	1	0
2	D	4	0	0	0	0
All	All	4119	0	3998	95	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 12.

All (95) 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:176:GLN:NE2	1:B:169:SER:OG	2.17	0.77
1:B:186:GLN:HE22	1:D:186:GLN:HG2	1.52	0.75
1:B:168:LEU:HB2	1:C:175:GLN:HE22	1.51	0.73
1:D:151:LEU:HG	1:C:190:MSE:HE2	1.70	0.73
1:A:164:GLN:HB3	1:D:178:ILE:HD13	1.72	0.70
1:B:132:GLN:HE21	1:B:133:GLN:HE21	1.43	0.67
1:B:164:GLN:HB2	1:C:179:PHE:HD1	1.60	0.66
1:B:175:GLN:HG3	1:C:168:LEU:HD13	1.78	0.66
1:A:197:HIS:HD2	1:B:148:MSE:HE2	1.62	0.65
1:A:124:LYS:HE2	1:A:128:ILE:HD11	1.77	0.64
1:B:185:VAL:HG11	1:C:157:ASP:OD2	1.97	0.63
1:B:132:GLN:NE2	1:B:133:GLN:HE21	1.96	0.63
1:B:186:GLN:NE2	1:D:186:GLN:HG2	2.14	0.63
1:D:144:SER:HA	1:C:197:HIS:CE1	2.35	0.62
1:D:155:GLU:HB2	1:C:190:MSE:HE1	1.81	0.62
1:B:197:HIS:HD2	1:C:143:TYR:OH	1.83	0.61
1:D:122:ASN:HD21	1:C:226:MSE:HG3	1.65	0.61
1:B:168:LEU:HB2	1:C:175:GLN:NE2	2.16	0.60
1:A:153:GLN:OE1	1:D:189:ARG:NH1	2.35	0.60
1:D:201:ILE:HD11	1:C:144:SER:HB2	1.83	0.59
1:D:166:GLU:O	1:D:169:SER:OG	2.22	0.57
1:D:169:SER:HB2	1:C:176:GLN:HE22	1.68	0.57
1:A:151:LEU:HD13	1:B:194:LYS:HG3	1.85	0.57
1:A:161:PHE:CE2	1:D:185:VAL:HG11	2.40	0.56
1:A:130:LYS:O	1:A:134:GLU:HG3	2.06	0.56
1:D:169:SER:CB	1:C:176:GLN:HE22	2.19	0.55
1:A:221:GLU:HB3	1:D:128:ILE:HD13	1.89	0.55
1:B:133:GLN:NE2	1:D:132:GLN:HE22	2.05	0.54
1:B:158:ILE:HG12	1:C:186:GLN:OE1	2.08	0.54
1:D:185:VAL:HA	1:D:188:GLN:HG2	1.90	0.54
1:A:162:GLU:HB2	1:B:183:ARG:NE	2.23	0.54
1:A:204:LEU:HD13	1:B:140:ASN:ND2	2.23	0.53
1:A:148:MSE:HE2	1:B:197:HIS:ND1	2.23	0.53
1:B:129:TRP:HD1	1:D:129:TRP:CD1	2.26	0.52
1:C:132:GLN:HE21	1:C:133:GLN:HG2	1.74	0.52
1:A:125:ILE:HD12	1:B:226:MSE:HE2	1.92	0.52
1:A:222:LEU:CD2	1:D:125:ILE:HG23	2.42	0.50
1:B:132:GLN:NE2	1:D:133:GLN:OE1	2.43	0.50
1:B:221:GLU:HB3	1:C:128:ILE:HD13	1.93	0.50
1:A:197:HIS:CD2	1:B:148:MSE:HE2	2.46	0.49
1:D:224:LYS:O	1:D:228:MSE:HG2	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:170:ASN:HA	1:A:173:ARG:HG2	1.94	0.49
1:A:118:PHE:CE1	1:D:229:LEU:HD21	2.48	0.48
1:A:162:GLU:HB2	1:B:183:ARG:HE	1.78	0.48
1:C:202:LYS:HE3	2:C:301:HOH:O	2.12	0.48
1:B:154:TRP:CD1	1:B:158:ILE:HG13	2.49	0.47
1:A:201:ILE:HD11	1:B:144:SER:HB2	1.95	0.47
1:B:147:PHE:CE1	1:D:147:PHE:HE1	2.33	0.47
1:D:188:GLN:HG3	1:D:189:ARG:N	2.29	0.47
1:B:133:GLN:CD	1:D:132:GLN:HE22	2.23	0.46
1:A:121:SER:O	1:A:125:ILE:HG13	2.16	0.46
1:A:161:PHE:HE2	1:D:185:VAL:HG11	1.81	0.46
1:D:142:GLU:O	1:D:146:GLN:HG3	2.15	0.46
1:C:167:LYS:O	1:C:171:LEU:HG	2.16	0.46
1:A:222:LEU:HD21	1:D:125:ILE:HG23	1.97	0.45
1:D:176:GLN:O	1:D:180:GLN:HG3	2.16	0.45
1:A:149:ASN:O	1:A:152:GLN:HG2	2.16	0.45
1:B:178:ILE:HD12	1:C:168:LEU:HD11	1.99	0.45
1:A:155:GLU:CD	1:B:190:MSE:HE1	2.42	0.45
1:B:230:GLN:O	1:B:234:MSE:HG3	2.16	0.45
1:B:202:LYS:HB2	1:B:202:LYS:HE2	1.71	0.45
1:B:213:ASN:HB3	1:B:217:GLY:N	2.32	0.45
1:A:169:SER:HB2	1:B:176:GLN:OE1	2.17	0.44
1:A:182:SER:O	1:A:186:GLN:HG2	2.17	0.44
1:B:129:TRP:HD1	1:D:129:TRP:HD1	1.65	0.44
1:B:138:LYS:HE2	1:B:142:GLU:OE1	2.18	0.44
1:C:167:LYS:H	1:C:167:LYS:HG2	1.59	0.44
1:D:173:ARG:HA	1:D:176:GLN:HG2	1.99	0.44
1:B:222:LEU:HD12	1:C:125:ILE:HG12	2.00	0.44
1:B:133:GLN:HE22	1:D:133:GLN:HE22	1.64	0.44
1:D:169:SER:HB2	1:C:176:GLN:NE2	2.31	0.44
1:D:172:PHE:CZ	1:C:172:PHE:HB2	2.53	0.43
1:A:129:TRP:CZ2	1:C:129:TRP:HB2	2.54	0.43
1:C:163:GLU:HG3	1:C:164:GLN:N	2.33	0.43
1:A:230:GLN:HE21	1:B:118:PHE:HB2	1.83	0.43
1:A:234:MSE:SE	1:B:114:THR:HG21	2.69	0.43
1:D:201:ILE:HD11	1:C:144:SER:CB	2.48	0.43
1:B:189:ARG:HD2	1:B:189:ARG:HA	1.89	0.42
1:C:163:GLU:O	1:C:167:LYS:HG2	2.18	0.42
1:A:124:LYS:O	1:A:128:ILE:HG12	2.20	0.42
1:A:190:MSE:SE	1:D:154:TRP:HH2	2.52	0.42
1:A:213:ASN:HD22	1:A:216:THR:H	1.66	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:144:SER:HA	1:C:197:HIS:HE1	1.81	0.42
1:D:172:PHE:CE2	1:C:172:PHE:CG	3.08	0.42
1:B:200:PHE:HB2	1:C:143:TYR:CD1	2.55	0.41
1:B:129:TRP:CD1	1:D:129:TRP:HD1	2.39	0.41
1:B:175:GLN:CG	1:C:168:LEU:HD13	2.50	0.41
1:A:156:LEU:HB3	1:A:160:LYS:HE2	2.02	0.41
1:D:133:GLN:HA	1:D:133:GLN:NE2	2.36	0.41
1:A:222:LEU:O	1:A:226:MSE:HG2	2.20	0.40
1:B:228:MSE:O	1:B:231:LYS:HG2	2.20	0.40
1:A:181:GLN:HA	1:A:184:ILE:HG22	2.03	0.40
1:C:189:ARG:O	1:C:193:ILE:HG23	2.21	0.40
1:D:192:ALA:O	1:D:196:ILE:HG13	2.20	0.40
1:C:231:LYS:HE3	1:C:231:LYS:HB2	1.88	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	126/144 (88%)	125 (99%)	0	1 (1%)	16	20
1	B	125/144 (87%)	124 (99%)	1 (1%)	0	100	100
1	C	121/144 (84%)	120 (99%)	1 (1%)	0	100	100
1	D	117/144 (81%)	116 (99%)	1 (1%)	0	100	100
All	All	489/576 (85%)	485 (99%)	3 (1%)	1 (0%)	43	55

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	217	GLY

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	116/128 (91%)	107 (92%)	9 (8%)	11	16
1	B	114/128 (89%)	104 (91%)	10 (9%)	9	12
1	C	108/128 (84%)	93 (86%)	15 (14%)	3	4
1	D	100/128 (78%)	91 (91%)	9 (9%)	9	12
All	All	438/512 (86%)	395 (90%)	43 (10%)	7	10

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	126	GLU
1	A	129	TRP
1	A	144	SER
1	A	166	GLU
1	A	169	SER
1	A	176	GLN
1	A	187	SER
1	A	194	LYS
1	A	214	LEU
1	B	132	GLN
1	B	141	ASN
1	B	148	MSE
1	B	177	LYS
1	B	182	SER
1	B	189	ARG
1	B	190	MSE
1	B	193	ILE
1	B	195	GLN
1	B	207	VAL
1	D	122	ASN
1	D	133	GLN
1	D	158	ILE
1	D	160	LYS
1	D	161	PHE

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Mol	Chain	Res	Type
1	D	190	MSE
1	D	207	VAL
1	D	219	GLN
1	D	233	VAL
1	C	142	GLU
1	C	150	VAL
1	C	159	GLN
1	C	163	GLU
1	C	175	GLN
1	C	184	ILE
1	C	185	VAL
1	C	186	GLN
1	C	191	LYS
1	C	193	ILE
1	C	207	VAL
1	C	214	LEU
1	C	221	GLU
1	C	228	MSE
1	C	234	MSE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	127	GLN
1	A	141	ASN
1	A	176	GLN
1	A	213	ASN
1	A	230	GLN
1	B	123	GLN
1	B	132	GLN
1	B	137	GLN
1	B	140	ASN
1	B	174	GLN
1	B	175	GLN
1	B	186	GLN
1	B	195	GLN
1	B	197	HIS
1	B	211	ASN
1	D	122	ASN
1	D	132	GLN
1	D	133	GLN
1	D	140	ASN

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Mol	Chain	Res	Type
1	D	145	GLN
1	D	152	GLN
1	D	153	GLN
1	D	159	GLN
1	D	164	GLN
1	D	174	GLN
1	D	175	GLN
1	D	176	GLN
1	D	181	GLN
1	C	123	GLN
1	C	132	GLN
1	C	133	GLN
1	C	140	ASN
1	C	145	GLN
1	C	146	GLN
1	C	153	GLN
1	C	159	GLN
1	C	176	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 ⓘ

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	121/144 (84%)	1.32	26 (21%) 2 3	44, 74, 94, 110	0
1	B	120/144 (83%)	1.25	25 (20%) 2 3	50, 73, 98, 128	0
1	C	118/144 (81%)	1.45	25 (21%) 2 3	45, 80, 103, 127	0
1	D	114/144 (79%)	1.60	32 (28%) 1 1	41, 75, 105, 123	0
All	All	473/576 (82%)	1.40	108 (22%) 2 2	41, 75, 101, 128	0

All (108) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	157	ASP	6.6
1	D	211	ASN	6.2
1	D	237	THR	6.0
1	C	222	LEU	5.2
1	C	214	LEU	4.9
1	B	237	THR	4.8
1	A	207	VAL	4.8
1	A	147	PHE	4.7
1	D	170	ASN	4.3
1	D	218	THR	4.2
1	C	212	ASP	4.2
1	B	168	LEU	4.2
1	C	207	VAL	4.1
1	A	180	GLN	4.0
1	C	237	THR	3.9
1	A	198	GLU	3.9
1	A	113	TYR	3.8
1	C	215	PHE	3.8
1	C	171	LEU	3.7
1	B	233	VAL	3.6
1	A	129	TRP	3.6

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Mol	Chain	Res	Type	RSRZ
1	D	222	LEU	3.4
1	C	161	PHE	3.3
1	C	127	GLN	3.3
1	A	171	LEU	3.2
1	A	136	ILE	3.2
1	A	187	SER	3.2
1	D	166	GLU	3.1
1	A	161	PHE	3.1
1	B	181	GLN	3.1
1	A	183	ARG	3.1
1	C	113	TYR	3.0
1	D	114	THR	3.0
1	A	116	ALA	3.0
1	D	171	LEU	2.9
1	D	199	GLN	2.9
1	D	178	ILE	2.9
1	D	129	TRP	2.9
1	A	222	LEU	2.9
1	C	164	GLN	2.9
1	D	197	HIS	2.9
1	D	169	SER	2.9
1	A	122	ASN	2.9
1	D	193	ILE	2.8
1	B	197	HIS	2.7
1	B	227	ALA	2.7
1	D	113	TYR	2.7
1	D	207	VAL	2.7
1	C	168	LEU	2.7
1	A	132	GLN	2.7
1	A	114	THR	2.6
1	A	193	ILE	2.6
1	C	151	LEU	2.6
1	B	179	PHE	2.6
1	C	129	TRP	2.6
1	C	137	GLN	2.5
1	B	236	GLU	2.5
1	D	179	PHE	2.5
1	D	233	VAL	2.5
1	C	132	GLN	2.5
1	C	117	SER	2.5
1	C	211	ASN	2.4
1	D	130	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	154	TRP	2.4
1	B	188	GLN	2.4
1	C	153	GLN	2.4
1	A	175	GLN	2.4
1	D	137	GLN	2.4
1	B	211	ASN	2.4
1	A	238	GLN	2.4
1	C	139	LEU	2.3
1	D	184	ILE	2.3
1	D	142	GLU	2.3
1	D	162	GLU	2.3
1	D	236	GLU	2.3
1	C	187	SER	2.3
1	B	175	GLN	2.3
1	B	113	TYR	2.2
1	B	114	THR	2.2
1	A	203	SER	2.2
1	D	182	SER	2.2
1	D	168	LEU	2.2
1	D	200	PHE	2.2
1	C	192	ALA	2.2
1	B	151	LEU	2.2
1	B	158	ILE	2.2
1	B	134	GLU	2.2
1	B	238	GLN	2.2
1	D	180	GLN	2.2
1	A	178	ILE	2.1
1	B	161	PHE	2.1
1	B	165	GLY	2.1
1	A	218	THR	2.1
1	B	150	VAL	2.1
1	A	117	SER	2.1
1	A	143	TYR	2.1
1	D	153	GLN	2.1
1	B	120	ALA	2.1
1	B	196	ILE	2.1
1	C	150	VAL	2.1
1	A	140	ASN	2.0
1	B	189	ARG	2.0
1	B	129	TRP	2.0
1	D	151	LEU	2.0
1	D	204	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	196	ILE	2.0
1	B	193	ILE	2.0
1	C	193	ILE	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.