



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:44 AM UTC

PDB ID : 7E5N / pdb_00007e5n
Title : crystal structure of cis assembled TROP-2
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Deposited on : 2021-02-19
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 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

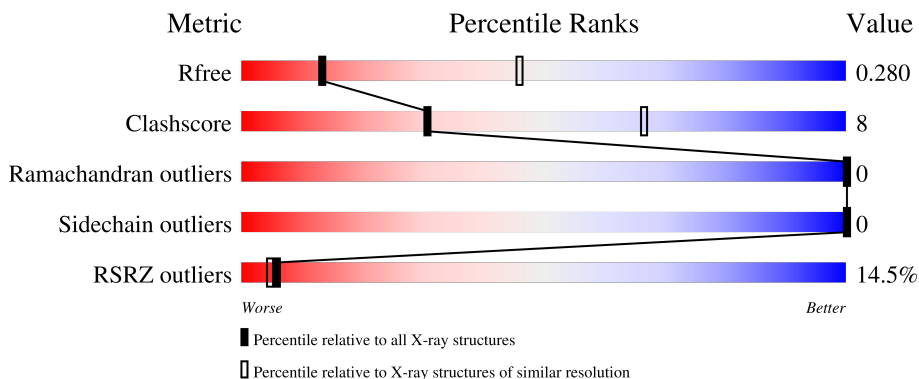
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	180053	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (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 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	323	8% 56% 17% • 27%
1	B	323	6% 61% 12% • 27%
1	C	323	9% 59% 12% • 29%
1	D	323	19% 56% 17% • 27%

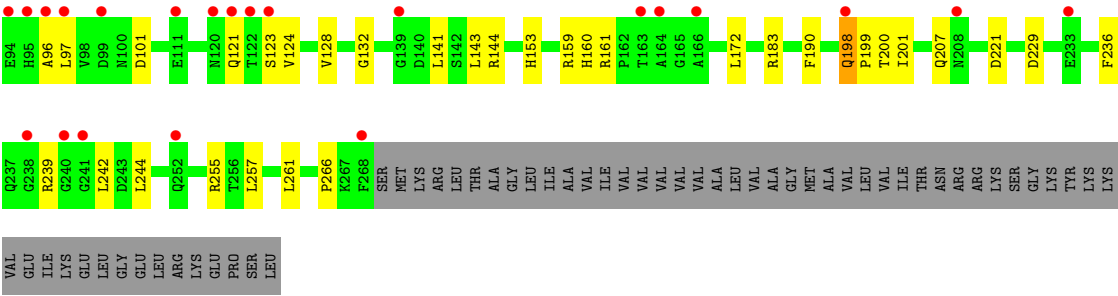
2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 7399 atoms, of which 0 are hydrogens and 0 are deuteriums.

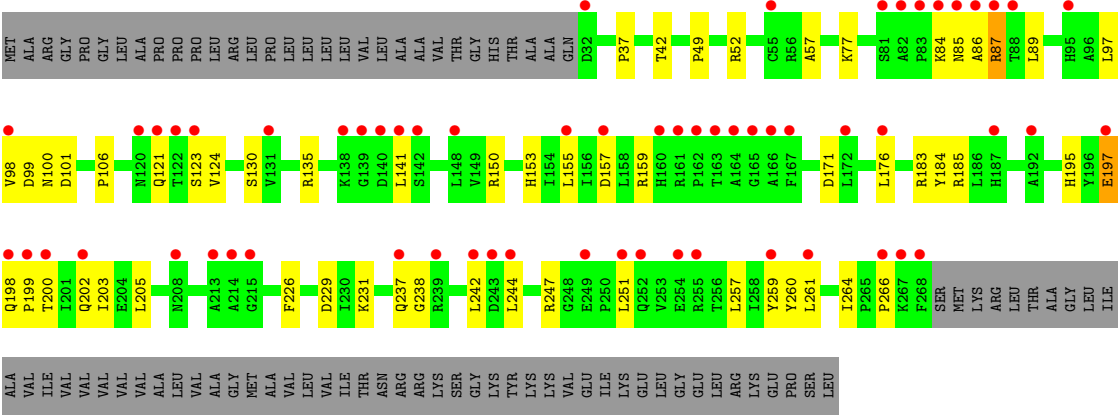
In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Tumor-associated calcium signal transducer 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	237	Total	C	N	O	S	0	0	0
			1864	1148	351	350	15			
1	B	237	Total	C	N	O	S	0	0	0
			1864	1148	351	350	15			
1	C	230	Total	C	N	O	S	0	0	0
			1807	1114	337	341	15			
1	D	237	Total	C	N	O	S	0	0	0
			1864	1148	351	350	15			



● Molecule 1: Tumor-associated calcium signal transducer 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 2 2	Depositor
Cell constants a, b, c, α , β , γ	145.14Å 145.14Å 217.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.95 – 3.20 49.95 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.5 (49.95-3.20) 99.5 (49.95-3.20)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.71 (at 3.19Å)	Xtriage
Refinement program	PHENIX (1.16_3549: ???)	Depositor
R, R_{free}	0.241 , 0.276 0.245 , 0.280	Depositor DCC
R_{free} test set	1946 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	70.9	Xtriage
Anisotropy	0.131	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 69.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	7399	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.12% 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	0.16	0/1899	0.48	0/2566
1	B	0.14	0/1899	0.62	5/2566 (0.2%)
1	C	0.13	0/1841	0.46	0/2485
1	D	0.16	0/1899	0.50	0/2566
All	All	0.15	0/7538	0.52	5/10183 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	B	0	4
1	C	0	1
1	D	0	3
All	All	0	11

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	121	GLN	CA-C-N	8.69	140.56	122.41
1	B	121	GLN	C-N-CA	8.69	140.56	122.41
1	B	120	ASN	CA-C-N	8.01	136.84	121.54
1	B	120	ASN	C-N-CA	8.01	136.84	121.54
1	B	122	THR	N-CA-C	-5.68	98.58	107.73

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	198	GLN	Peptide
1	A	80	MET	Peptide
1	A	87	ARG	Peptide
1	B	120	ASN	Peptide
1	B	121	GLN	Peptide
1	B	198	GLN	Peptide
1	B	85	ASN	Peptide
1	C	198	GLN	Peptide
1	D	197	GLU	Peptide
1	D	198	GLN	Peptide
1	D	87	ARG	Peptide

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	1864	0	1824	39	0
1	B	1864	0	1824	27	0
1	C	1807	0	1760	31	0
1	D	1864	0	1824	38	0
All	All	7399	0	7232	122	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 8.

All (122) 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:C:242:LEU:HD22	1:C:244:LEU:HB2	1.45	0.99
1:A:86:ALA:HB1	1:A:89:LEU:HB2	1.44	0.98
1:D:86:ALA:HB1	1:D:89:LEU:HB2	1.54	0.89
1:C:229:ASP:HB3	1:C:242:LEU:HD12	1.63	0.79
1:A:183:ARG:O	1:A:185:ARG:NH2	2.16	0.78
1:D:195:HIS:HB2	1:D:202:GLN:HB2	1.68	0.76
1:D:242:LEU:HD13	1:D:244:LEU:HB2	1.68	0.75
1:D:229:ASP:HB3	1:D:242:LEU:HB2	1.69	0.74
1:A:242:LEU:HD22	1:A:244:LEU:HB2	1.68	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:97:LEU:HD21	1:D:266:PRO:HG3	1.68	0.73
1:A:37:PRO:HA	1:A:237:GLN:HB2	1.71	0.72
1:C:42:THR:HG22	1:C:57:ALA:HA	1.70	0.71
1:B:93:SER:HB3	1:B:96:ALA:HB2	1.74	0.69
1:C:92:PRO:O	1:D:202:GLN:NE2	2.27	0.67
1:B:77:LYS:NZ	1:B:106:PRO:O	2.28	0.66
1:A:85:ASN:HB2	1:A:87:ARG:HG3	1.78	0.65
1:D:77:LYS:NZ	1:D:106:PRO:O	2.30	0.65
1:D:242:LEU:CD1	1:D:244:LEU:HB2	2.27	0.64
1:C:183:ARG:NH2	1:C:229:ASP:OD2	2.31	0.63
1:B:175:GLU:OE1	1:B:178:ARG:NH2	2.32	0.62
1:D:183:ARG:O	1:D:185:ARG:NH2	2.32	0.62
1:C:159:ARG:HG2	1:C:200:THR:HG22	1.80	0.61
1:D:197:GLU:O	1:D:200:THR:N	2.35	0.59
1:C:153:HIS:HB3	1:C:261:LEU:HB3	1.86	0.58
1:B:121:GLN:O	1:B:121:GLN:HG2	2.04	0.57
1:B:169:HIS:CD2	1:B:196:TYR:HB3	2.40	0.57
1:C:229:ASP:CB	1:C:242:LEU:HD12	2.36	0.55
1:B:183:ARG:O	1:B:185:ARG:NH2	2.40	0.55
1:A:255:ARG:HG2	1:A:256:THR:H	1.72	0.55
1:D:155:LEU:HB3	1:D:259:TYR:HB2	1.88	0.54
1:D:176:LEU:HD13	1:D:203:ILE:HD11	1.89	0.54
1:A:167:PHE:HE1	1:A:249:GLU:HG2	1.73	0.54
1:D:37:PRO:HA	1:D:237:GLN:HB2	1.90	0.54
1:A:183:ARG:NH2	1:A:229:ASP:OD2	2.33	0.54
1:C:255:ARG:HH21	1:C:257:LEU:HD12	1.72	0.53
1:D:171:ASP:OD2	1:D:247:ARG:NH1	2.41	0.53
1:B:33:ASN:OD1	1:B:34:CYS:N	2.42	0.53
1:B:94:GLU:HG3	1:B:95:HIS:ND1	2.23	0.53
1:A:123:SER:HA	1:A:141:LEU:HD21	1.90	0.53
1:C:121:GLN:HB3	1:C:124:VAL:HG22	1.91	0.52
1:A:109:ASP:HB2	1:A:110:PRO:HD2	1.92	0.52
1:A:155:LEU:HD23	1:B:92:PRO:HA	1.91	0.52
1:D:135:ARG:NH2	1:D:264:ILE:O	2.44	0.51
1:D:49:PRO:HG2	1:D:52:ARG:HH21	1.75	0.51
1:D:85:ASN:OD1	1:D:87:ARG:NE	2.38	0.51
1:B:150:ARG:NH1	1:B:218:ASP:OD2	2.33	0.51
1:D:98:VAL:HG22	1:D:99:ASP:H	1.76	0.51
1:A:57:ALA:HB2	1:A:64:VAL:HG13	1.91	0.51
1:C:101:ASP:OD1	1:D:231:LYS:NZ	2.30	0.50
1:A:227:GLU:HG3	1:A:228:ARG:H	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:179:LEU:HD11	1:A:242:LEU:HD21	1.94	0.50
1:A:266:PRO:HG3	1:B:97:LEU:HD11	1.94	0.49
1:D:101:ASP:OD2	1:D:101:ASP:N	2.44	0.49
1:D:244:LEU:HG	1:D:251:LEU:HB3	1.94	0.49
1:B:160:HIS:CE1	1:B:199:PRO:HA	2.48	0.49
1:C:255:ARG:NH2	1:C:257:LEU:HD12	2.28	0.48
1:D:183:ARG:NH2	1:D:242:LEU:HD23	2.28	0.48
1:C:159:ARG:HG3	1:C:255:ARG:NH1	2.28	0.48
1:A:159:ARG:HD3	1:A:255:ARG:NH2	2.29	0.47
1:B:42:THR:HG22	1:B:57:ALA:HA	1.96	0.47
1:C:143:LEU:HD23	1:C:144:ARG:N	2.28	0.47
1:C:160:HIS:CE1	1:C:199:PRO:HA	2.50	0.47
1:B:57:ALA:HB2	1:B:64:VAL:HG13	1.96	0.47
1:B:229:ASP:HB3	1:B:242:LEU:HG	1.95	0.47
1:A:78:ALA:O	1:A:81:SER:HB3	2.14	0.47
1:D:153:HIS:ND1	1:D:261:LEU:HD13	2.30	0.47
1:A:158:LEU:HD22	1:A:253:VAL:HG11	1.97	0.47
1:A:227:GLU:HG3	1:A:228:ARG:N	2.31	0.46
1:C:236:PHE:CD1	1:C:239:ARG:HD3	2.50	0.46
1:D:98:VAL:HG13	1:D:100:ASN:H	1.80	0.46
1:C:101:ASP:HA	1:D:100:ASN:ND2	2.30	0.45
1:A:85:ASN:HB2	1:A:87:ARG:CG	2.47	0.45
1:A:130:SER:HA	1:A:148:LEU:HD11	1.97	0.45
1:C:128:VAL:HB	1:C:132:GLY:HA2	1.98	0.45
1:B:190:PHE:CE1	1:B:207:GLN:HG3	2.52	0.45
1:A:255:ARG:HG2	1:A:256:THR:N	2.31	0.45
1:A:131:VAL:HG13	1:A:133:VAL:HG22	1.99	0.44
1:A:183:ARG:HH22	1:A:242:LEU:HA	1.82	0.44
1:B:84:LYS:H	1:B:84:LYS:HG3	1.64	0.44
1:A:35:THR:O	1:C:52:ARG:NH2	2.50	0.44
1:C:161:ARG:NH2	1:D:238:GLY:O	2.51	0.44
1:D:121:GLN:HB3	1:D:124:VAL:HB	2.00	0.44
1:A:244:LEU:HG	1:A:251:LEU:HB3	1.99	0.44
1:C:266:PRO:HG3	1:D:97:LEU:HD11	1.98	0.44
1:D:42:THR:HG22	1:D:57:ALA:HA	2.00	0.43
1:A:157:ASP:HB3	1:A:257:LEU:HB2	1.99	0.43
1:B:104:TYR:O	1:B:106:PRO:HD3	2.18	0.43
1:A:125:CYS:SG	1:A:141:LEU:HD23	2.59	0.43
1:B:198:GLN:HG3	1:B:199:PRO:HD3	1.99	0.43
1:D:157:ASP:HB3	1:D:257:LEU:HB2	2.01	0.43
1:C:72:LYS:N	1:C:221:ASP:OD1	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:198:GLN:O	1:A:200:THR:N	2.52	0.42
1:A:52:ARG:HH11	1:C:53:CYS:HB3	1.85	0.42
1:B:150:ARG:NH2	1:B:207:GLN:OE1	2.49	0.42
1:B:130:SER:HA	1:B:148:LEU:HD11	2.01	0.42
1:B:242:LEU:HA	1:B:242:LEU:HD23	1.80	0.42
1:A:134:ARG:HD2	1:A:136:THR:O	2.20	0.42
1:D:130:SER:HB3	1:D:150:ARG:HA	2.00	0.42
1:A:172:LEU:HD21	1:A:201:ILE:HD13	2.02	0.42
1:A:255:ARG:HG3	1:B:80:MET:HE3	2.01	0.42
1:C:172:LEU:HD21	1:C:201:ILE:HD13	2.02	0.42
1:C:198:GLN:HG3	1:C:199:PRO:HD3	2.01	0.42
1:D:153:HIS:HA	1:D:205:LEU:O	2.19	0.42
1:C:96:ALA:HB1	1:D:260:TYR:O	2.20	0.41
1:D:159:ARG:HG3	1:D:199:PRO:HG2	2.02	0.41
1:B:196:TYR:CD1	1:B:201:ILE:HG12	2.55	0.41
1:A:227:GLU:HB3	1:A:258:ILE:HD11	2.02	0.41
1:D:184:TYR:CZ	1:D:226:PHE:HA	2.56	0.41
1:B:261:LEU:HD23	1:B:261:LEU:HA	1.91	0.41
1:C:97:LEU:O	1:D:260:TYR:N	2.40	0.41
1:A:159:ARG:NH2	1:A:254:GLU:OE1	2.54	0.41
1:C:123:SER:HA	1:C:141:LEU:HG	2.02	0.41
1:C:190:PHE:CE1	1:C:207:GLN:HG3	2.56	0.41
1:A:84:LYS:HG2	1:A:85:ASN:H	1.85	0.41
1:A:121:GLN:HB3	1:A:124:VAL:HB	2.03	0.41
1:B:40:LYS:HA	1:B:237:GLN:HE21	1.86	0.41
1:B:57:ALA:HB2	1:B:64:VAL:CG1	2.49	0.41
1:D:123:SER:HA	1:D:141:LEU:HD21	2.03	0.41
1:A:102:GLY:O	1:A:104:TYR:HD1	2.03	0.40
1:A:176:LEU:HD13	1:A:203:ILE:HD11	2.03	0.40
1:C:37:PRO:HD2	1:C:66:CYS:HB3	2.03	0.40
1:D:84:LYS:HG3	1:D:85:ASN:HD22	1.86	0.40

There are no symmetry-related clashes.

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	235/323 (73%)	232 (99%)	3 (1%)	0	100	100
1	B	235/323 (73%)	231 (98%)	4 (2%)	0	100	100
1	C	226/323 (70%)	224 (99%)	2 (1%)	0	100	100
1	D	235/323 (73%)	229 (97%)	6 (3%)	0	100	100
All	All	931/1292 (72%)	916 (98%)	15 (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	204/272 (75%)	204 (100%)	0	100	100
1	B	204/272 (75%)	204 (100%)	0	100	100
1	C	198/272 (73%)	198 (100%)	0	100	100
1	D	204/272 (75%)	204 (100%)	0	100	100
All	All	810/1088 (74%)	810 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	160	HIS
1	B	120	ASN
1	B	169	HIS
1	B	237	GLN
1	C	211	GLN
1	D	211	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	237/323 (73%)	0.58	26 (10%) 10 7	29, 58, 130, 185	0
1	B	237/323 (73%)	0.65	20 (8%) 17 11	37, 65, 131, 198	0
1	C	230/323 (71%)	0.67	29 (12%) 8 6	37, 65, 144, 172	0
1	D	237/323 (73%)	1.28	61 (25%) 1 2	53, 110, 166, 200	0
All	All	941/1292 (72%)	0.79	136 (14%) 6 5	29, 72, 150, 200	0

All (136) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	92	PRO	7.4
1	D	200	THR	5.5
1	C	83	PRO	5.4
1	A	122	THR	5.2
1	A	82	ALA	5.1
1	C	94	GLU	5.1
1	D	163	THR	4.9
1	D	167	PHE	4.8
1	B	122	THR	4.7
1	B	123	SER	4.6
1	D	164	ALA	4.6
1	A	123	SER	4.6
1	C	96	ALA	4.5
1	A	86	ALA	4.5
1	D	161	ARG	4.4
1	C	84	LYS	4.3
1	C	97	LEU	4.3
1	B	198	GLN	4.1
1	A	166	ALA	4.1
1	B	164	ALA	4.0
1	A	81	SER	4.0

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Mol	Chain	Res	Type	RSRZ
1	B	121	GLN	4.0
1	B	208	ASN	3.9
1	B	240	GLY	3.9
1	D	86	ALA	3.9
1	D	254	GLU	3.8
1	A	84	LYS	3.7
1	B	120	ASN	3.7
1	D	123	SER	3.6
1	D	267	LYS	3.6
1	B	85	ASN	3.6
1	D	239	ARG	3.5
1	C	122	THR	3.5
1	A	198	GLN	3.5
1	D	202	GLN	3.5
1	C	166	ALA	3.4
1	A	199	PRO	3.4
1	B	83	PRO	3.4
1	D	199	PRO	3.3
1	B	86	ALA	3.3
1	C	238	GLY	3.3
1	A	121	GLN	3.3
1	D	157	ASP	3.3
1	C	240	GLY	3.2
1	D	166	ALA	3.2
1	C	99	ASP	3.1
1	A	144	ARG	3.1
1	D	259	TYR	3.1
1	C	164	ALA	3.1
1	D	162	PRO	3.1
1	B	163	THR	3.1
1	D	208	ASN	3.0
1	C	121	GLN	3.0
1	B	84	LYS	3.0
1	D	172	LEU	3.0
1	D	197	GLU	3.0
1	D	176	LEU	3.0
1	D	243	ASP	3.0
1	D	120	ASN	3.0
1	D	252	GLN	2.9
1	D	88	THR	2.9
1	D	198	GLN	2.9
1	D	261	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	33	ASN	2.9
1	D	81	SER	2.8
1	D	142	SER	2.8
1	C	95	HIS	2.8
1	C	120	ASN	2.8
1	D	85	ASN	2.8
1	D	82	ALA	2.8
1	D	84	LYS	2.8
1	A	211	GLN	2.8
1	C	32	ASP	2.8
1	D	87	ARG	2.8
1	C	252	GLN	2.8
1	D	215	GLY	2.7
1	C	208	ASN	2.7
1	A	80	MET	2.7
1	A	210	SER	2.7
1	B	92	PRO	2.7
1	D	187	HIS	2.6
1	A	268	PHE	2.6
1	B	165	GLY	2.6
1	D	192	ALA	2.6
1	D	148	LEU	2.6
1	D	255	ARG	2.6
1	A	35	THR	2.5
1	D	122	THR	2.5
1	A	87	ARG	2.5
1	D	138	LYS	2.5
1	D	213	ALA	2.4
1	D	214	ALA	2.4
1	D	268	PHE	2.4
1	A	83	PRO	2.4
1	C	139	GLY	2.4
1	A	165	GLY	2.4
1	D	165	GLY	2.4
1	D	121	GLN	2.4
1	A	32	ASP	2.3
1	D	140	ASP	2.3
1	A	251	LEU	2.3
1	C	123	SER	2.3
1	B	144	ARG	2.3
1	C	241	GLY	2.3
1	A	163	THR	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	209	THR	2.3
1	D	244	LEU	2.3
1	C	198	GLN	2.3
1	D	139	GLY	2.3
1	B	81	SER	2.3
1	D	251	LEU	2.2
1	B	241	GLY	2.2
1	D	131	VAL	2.2
1	B	247	ARG	2.2
1	D	160	HIS	2.2
1	C	81	SER	2.2
1	D	237	GLN	2.2
1	B	88	THR	2.2
1	D	141	LEU	2.2
1	A	208	ASN	2.2
1	A	255	ARG	2.2
1	D	98	VAL	2.1
1	C	268	PHE	2.1
1	D	55	CYS	2.1
1	C	111	GLU	2.1
1	D	249	GLU	2.1
1	D	155	LEU	2.1
1	D	83	PRO	2.1
1	D	32	ASP	2.1
1	C	163	THR	2.1
1	C	93	SER	2.1
1	C	233	GLU	2.1
1	D	266	PRO	2.1
1	A	85	ASN	2.0
1	D	242	LEU	2.0
1	D	95	HIS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

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.