



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 10, 2023 – 06:04 AM EDT

PDB ID : 7SD1
Title : Crystal structure of SHOC2
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Deposited on : 2021-09-29
Resolution : 3.19 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.35.1
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) : 2.35.1

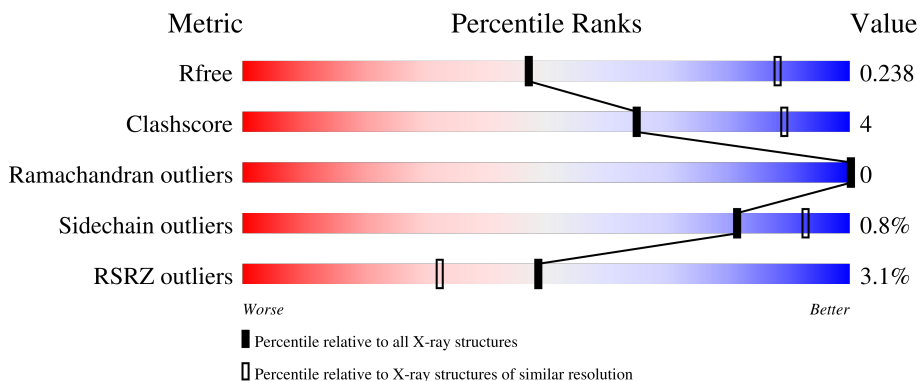
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.19 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 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	585	 3% 77% 8% 12%
1	B	585	 2% 73% 12% 13%
1	C	585	 2% 75% 10% 13%
1	D	585	 3% 75% 11% 11%

2 Entry composition i

There is only 1 type of molecule in this entry. The entry contains 15721 atoms, of which 0 are hydrogens and 0 are deuteriums.

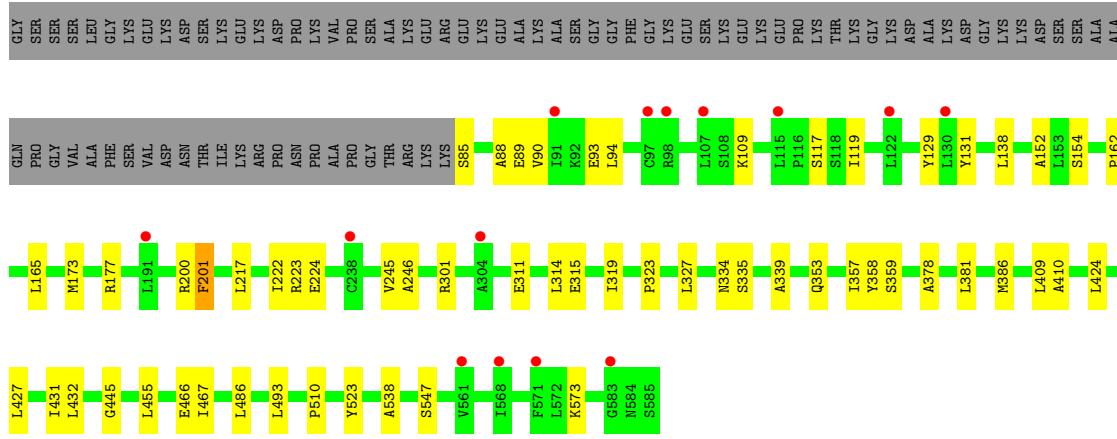
In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Leucine-rich repeat protein SHOC-2.

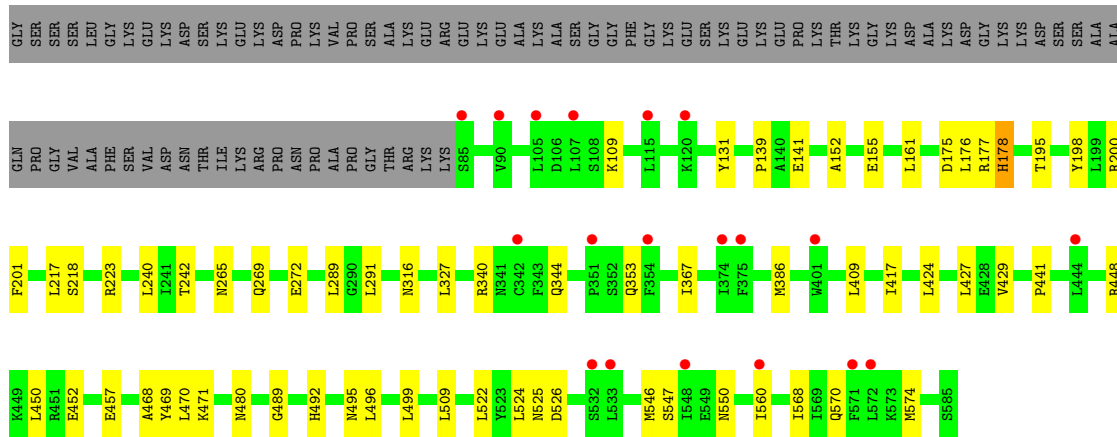
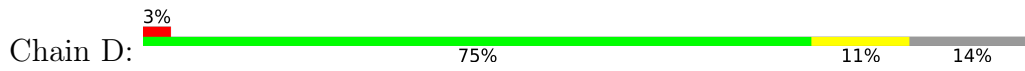
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	497	3916	2478	674	746	18	0	0	0
1	B	498	3923	2483	675	747	18	0	0	0
1	C	501	3941	2492	679	752	18	0	0	0
1	D	501	3941	2492	679	752	18	0	0	0

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP Q9UQ13
A	583	GLY	-	expression tag	UNP Q9UQ13
A	584	ASN	-	expression tag	UNP Q9UQ13
A	585	SER	-	expression tag	UNP Q9UQ13
B	1	GLY	-	expression tag	UNP Q9UQ13
B	583	GLY	-	expression tag	UNP Q9UQ13
B	584	ASN	-	expression tag	UNP Q9UQ13
B	585	SER	-	expression tag	UNP Q9UQ13
C	1	GLY	-	expression tag	UNP Q9UQ13
C	583	GLY	-	expression tag	UNP Q9UQ13
C	584	ASN	-	expression tag	UNP Q9UQ13
C	585	SER	-	expression tag	UNP Q9UQ13
D	1	GLY	-	expression tag	UNP Q9UQ13
D	583	GLY	-	expression tag	UNP Q9UQ13
D	584	ASN	-	expression tag	UNP Q9UQ13
D	585	SER	-	expression tag	UNP Q9UQ13



● Molecule 1: Leucine-rich repeat protein SHOC-2



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	168.63Å 201.83Å 233.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.13 – 3.19 49.13 – 3.19	Depositor EDS
% Data completeness (in resolution range)	96.3 (49.13-3.19) 82.8 (49.13-3.19)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.21 (at 3.19Å)	Xtrriage
Refinement program	PHENIX 1.19.1-4122_final	Depositor
R, R_{free}	0.208 , 0.238 0.209 , 0.238	Depositor DCC
R_{free} test set	1986 reflections (3.00%)	wwPDB-VP
Wilson B-factor (Å ²)	86.9	Xtrriage
Anisotropy	0.289	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 42.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	15721	wwPDB-VP
Average B, all atoms (Å ²)	120.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.88% 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	0.31	0/3973	0.55	0/5384
1	B	0.33	0/3980	0.54	0/5394
1	C	0.31	0/3998	0.53	0/5418
1	D	0.29	0/3998	0.53	0/5418
All	All	0.31	0/15949	0.54	0/21614

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	3916	0	4078	29	0
1	B	3923	0	4087	36	0
1	C	3941	0	4101	31	0
1	D	3941	0	4101	36	0
All	All	15721	0	16367	131	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 4.

All (131) 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:362:MET:HE1	1:A:367:ILE:HD13	1.74	0.69
1:B:85:SER:O	1:B:89:GLU:HG3	2.00	0.62
1:D:177:ARG:HG2	1:D:200:ARG:H	1.67	0.60
1:B:525:ASN:HB3	1:B:547:SER:HB3	1.83	0.59
1:A:362:MET:CE	1:A:367:ILE:CD1	2.81	0.59
1:D:522:LEU:O	1:D:546:MET:HA	2.04	0.58
1:B:149:MET:HA	1:B:171:LEU:HA	1.84	0.57
1:C:314:LEU:HD13	1:C:319:ILE:HD13	1.87	0.57
1:B:166:ASP:HB2	1:B:190:ARG:HH21	1.69	0.57
1:D:468:ALA:HB1	1:D:492:HIS:ND1	2.20	0.57
1:B:334:ASN:HA	1:B:357:ILE:HA	1.88	0.56
1:C:311:GLU:HG3	1:C:335:SER:HB3	1.89	0.55
1:C:386:MET:HB2	1:C:409:LEU:HD23	1.88	0.55
1:B:184:ILE:HD11	1:B:207:VAL:HG22	1.89	0.55
1:A:362:MET:HE2	1:A:367:ILE:CD1	2.36	0.55
1:A:362:MET:HE2	1:A:367:ILE:HD11	1.87	0.55
1:C:301:ARG:HG2	1:C:323:PRO:HB3	1.89	0.54
1:A:386:MET:HB2	1:A:409:LEU:HD23	1.90	0.54
1:C:201:PHE:H	1:C:224:GLU:HB2	1.72	0.54
1:B:579:ARG:HH21	1:B:581:MET:HA	1.71	0.54
1:A:218:SER:HA	1:A:240:LEU:HA	1.89	0.54
1:A:177:ARG:HG2	1:A:200:ARG:H	1.73	0.53
1:B:226:LYS:HG2	1:D:272:GLU:OE1	2.09	0.53
1:C:315:GLU:HG3	1:C:339:ALA:HB3	1.89	0.53
1:C:162:PRO:HG2	1:C:165:LEU:HD13	1.91	0.53
1:D:471:LYS:NZ	1:D:492:HIS:HB3	2.24	0.53
1:B:327:LEU:HB2	1:B:353:GLN:HB2	1.90	0.52
1:C:129:TYR:HA	1:C:152:ALA:HB3	1.92	0.52
1:B:467:ILE:O	1:B:493:LEU:HD21	2.09	0.52
1:A:424:LEU:HB3	1:A:427:LEU:HB2	1.92	0.51
1:D:327:LEU:HB2	1:D:353:GLN:HB2	1.92	0.51
1:B:223:ARG:HG2	1:B:224:GLU:HG2	1.93	0.51
1:A:362:MET:HE1	1:A:367:ILE:CD1	2.41	0.51
1:D:109:LYS:HE3	1:D:131:TYR:HE2	1.76	0.51
1:B:181:LEU:HD22	1:B:185:PRO:HG3	1.92	0.50
1:B:386:MET:HB2	1:B:409:LEU:HD23	1.93	0.50
1:C:523:TYR:HA	1:C:547:SER:HB2	1.95	0.49
1:C:357:ILE:HG22	1:C:359:SER:H	1.78	0.49
1:A:109:LYS:HE3	1:A:131:TYR:HE1	1.77	0.48
1:B:115:LEU:O	1:B:139:PRO:HG3	2.13	0.48
1:C:467:ILE:O	1:C:493:LEU:HD21	2.13	0.48
1:C:327:LEU:HB2	1:C:353:GLN:HB2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:424:LEU:HD23	1:D:427:LEU:HD13	1.95	0.48
1:B:311:GLU:HG3	1:B:335:SER:HB3	1.95	0.48
1:D:242:THR:HG23	1:D:265:ASN:HB3	1.95	0.48
1:D:218:SER:HA	1:D:240:LEU:HA	1.95	0.47
1:B:444:LEU:O	1:B:470:LEU:HD21	2.14	0.47
1:D:448:ARG:CZ	1:D:469:TYR:HB3	2.45	0.47
1:D:155:GLU:HA	1:D:178:HIS:HB3	1.96	0.46
1:A:255:LYS:HE3	1:A:279:THR:HG23	1.96	0.46
1:D:525:ASN:HB3	1:D:547:SER:HB3	1.96	0.46
1:B:218:SER:HA	1:B:240:LEU:HA	1.97	0.46
1:A:323:PRO:HD2	1:A:326:LEU:HD22	1.97	0.46
1:B:162:PRO:HG2	1:B:165:LEU:CD1	2.46	0.46
1:C:486:LEU:HB2	1:C:510:PRO:HD3	1.97	0.46
1:A:85:SER:HB2	1:A:110:ARG:O	2.16	0.45
1:A:424:LEU:HD23	1:A:427:LEU:HD13	1.98	0.45
1:A:334:ASN:HA	1:A:357:ILE:HA	1.98	0.45
1:C:200:ARG:HG3	1:C:223:ARG:HB3	1.98	0.45
1:D:450:LEU:HD23	1:D:470:LEU:HD13	1.99	0.45
1:A:327:LEU:HB2	1:A:353:GLN:HB2	1.98	0.45
1:C:424:LEU:HD23	1:C:427:LEU:HD13	1.98	0.45
1:D:152:ALA:HA	1:D:175:ASP:HB3	1.99	0.45
1:B:222:ILE:HD11	1:B:245:VAL:HG12	1.99	0.45
1:D:161:LEU:HD11	1:D:176:LEU:HD11	1.99	0.45
1:C:432:LEU:HB2	1:C:455:LEU:HD23	1.98	0.44
1:D:195:THR:HA	1:D:217:LEU:HA	1.99	0.44
1:C:89:GLU:O	1:C:93:GLU:HG2	2.17	0.44
1:A:315:GLU:HG3	1:A:339:ALA:HB3	1.99	0.44
1:B:526:ASP:HA	1:B:550:ASN:HB2	1.98	0.44
1:D:429:VAL:HG13	1:D:452:GLU:HB3	1.98	0.44
1:A:574:MET:O	1:A:579:ARG:HD3	2.18	0.44
1:B:196:THR:HG23	1:B:219:MET:HB3	1.99	0.44
1:D:509:LEU:HD11	1:D:524:LEU:HD11	2.00	0.44
1:B:181:LEU:HD12	1:B:202:ASN:ND2	2.33	0.43
1:C:424:LEU:HB3	1:C:427:LEU:HB2	2.00	0.43
1:D:417:ILE:HB	1:D:441:PRO:HD3	2.00	0.43
1:D:560:ILE:HG22	1:D:568:ILE:HG12	1.98	0.43
1:A:373:GLY:H	1:A:397:ASP:CG	2.22	0.43
1:D:344:GLN:O	1:D:367:ILE:HA	2.17	0.43
1:D:139:PRO:HB2	1:D:141:GLU:OE1	2.18	0.43
1:A:139:PRO:HB2	1:A:141:GLU:OE1	2.18	0.43
1:D:496:LEU:HD21	1:D:499:LEU:HD13	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:152:ALA:HA	1:A:175:ASP:HB3	1.99	0.43
1:B:242:THR:HG23	1:B:265:ASN:HB2	2.01	0.43
1:C:138:LEU:HD23	1:C:138:LEU:HA	1.78	0.43
1:D:457:GLU:HG2	1:D:480:ASN:ND2	2.33	0.43
1:C:200:ARG:HE	1:C:224:GLU:HG3	1.83	0.43
1:C:538:ALA:HB1	1:C:573:LYS:HG3	2.00	0.43
1:D:448:ARG:HA	1:D:448:ARG:HD3	1.73	0.43
1:A:181:LEU:O	1:A:203:ARG:N	2.50	0.43
1:B:440:LEU:HD11	1:B:455:LEU:CD1	2.49	0.43
1:C:223:ARG:HG3	1:C:246:ALA:HB3	2.01	0.43
1:A:118:SER:C	1:A:120:LYS:N	2.73	0.42
1:A:440:LEU:HD13	1:A:444:LEU:HD22	2.00	0.42
1:D:316:ASN:H	1:D:340:ARG:HB2	1.84	0.42
1:D:177:ARG:HG3	1:D:198:TYR:HB3	2.01	0.42
1:D:526:ASP:HA	1:D:550:ASN:HB2	2.01	0.42
1:D:386:MET:HB2	1:D:409:LEU:HD23	2.02	0.42
1:A:549:GLU:OE1	1:A:549:GLU:N	2.45	0.42
1:C:109:LYS:HE3	1:C:131:TYR:HE2	1.84	0.42
1:A:532:SER:HB3	1:A:555:HIS:HB2	2.02	0.42
1:B:539:LEU:HD21	1:B:578:TYR:HB3	2.00	0.42
1:C:334:ASN:HB2	1:C:358:TYR:HD1	1.85	0.42
1:C:378:ALA:HB1	1:C:381:LEU:HB2	2.01	0.42
1:A:522:LEU:O	1:A:546:MET:HA	2.20	0.42
1:B:122:LEU:O	1:B:145:LEU:HD22	2.20	0.42
1:B:344:GLN:HG2	1:B:366:ARG:HB3	2.02	0.42
1:C:117:SER:C	1:C:119:ILE:H	2.22	0.42
1:A:319:ILE:H	1:A:319:ILE:HG13	1.72	0.41
1:B:557:PRO:HA	1:B:558:PRO:HD3	1.95	0.41
1:D:471:LYS:O	1:D:495:ASN:ND2	2.50	0.41
1:B:212:LYS:HD3	1:B:233:GLU:HG2	2.01	0.41
1:B:334:ASN:HB2	1:B:358:TYR:HD1	1.85	0.41
1:C:445:GLY:HA3	1:C:466:GLU:HB3	2.02	0.41
1:D:269:GLN:H	1:D:269:GLN:HG2	1.71	0.41
1:D:289:LEU:HG	1:D:291:LEU:HG	2.02	0.41
1:B:522:LEU:O	1:B:546:MET:HA	2.20	0.41
1:B:138:LEU:HD23	1:B:138:LEU:HA	1.77	0.41
1:D:570:GLN:O	1:D:574:MET:HG2	2.21	0.41
1:C:177:ARG:HG2	1:C:200:ARG:H	1.86	0.41
1:C:410:ALA:HB2	1:C:431:ILE:HG22	2.02	0.41
1:D:468:ALA:HB2	1:D:489:GLY:HA3	2.03	0.41
1:B:490:ILE:O	1:B:516:LEU:HD21	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:90:VAL:O	1:C:94:LEU:HG	2.20	0.40
1:A:200:ARG:HG3	1:A:223:ARG:HB3	2.02	0.40
1:B:533:LEU:HD13	1:B:546:MET:HE1	2.04	0.40
1:B:138:LEU:O	1:B:162:PRO:HG3	2.22	0.40
1:C:85:SER:O	1:C:88:ALA:N	2.52	0.40
1:D:200:ARG:HG2	1:D:223:ARG:HB3	2.03	0.40
1:B:185:PRO:HB2	1:B:188:VAL:HG23	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	495/585 (85%)	477 (96%)	18 (4%)	0	100	100
1	B	496/585 (85%)	478 (96%)	18 (4%)	0	100	100
1	C	499/585 (85%)	485 (97%)	14 (3%)	0	100	100
1	D	499/585 (85%)	472 (95%)	27 (5%)	0	100	100
All	All	1989/2340 (85%)	1912 (96%)	77 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	462/532 (87%)	461 (100%)	1 (0%)	93	98
1	B	463/532 (87%)	458 (99%)	5 (1%)	73	88
1	C	465/532 (87%)	459 (99%)	6 (1%)	69	87
1	D	465/532 (87%)	463 (100%)	2 (0%)	91	95
All	All	1855/2128 (87%)	1841 (99%)	14 (1%)	81	93

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

Mol	Chain	Res	Type
1	A	201	PHE
1	B	154	SER
1	B	159	THR
1	B	210	ASP
1	B	448	ARG
1	B	508	HIS
1	C	154	SER
1	C	173	MET
1	C	201	PHE
1	C	217	LEU
1	C	222	ILE
1	C	245	VAL
1	D	178	HIS
1	D	201	PHE

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	247	HIS
1	A	316	ASN
1	A	344	GLN
1	A	364	HIS
1	C	446	ASN
1	D	178	HIS

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 monosaccharides 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	497/585 (84%)	0.11	15 (3%) 50 34	63, 106, 186, 250	0
1	B	498/585 (85%)	0.17	14 (2%) 53 37	69, 115, 184, 221	0
1	C	501/585 (85%)	0.17	14 (2%) 53 37	71, 112, 170, 215	0
1	D	501/585 (85%)	0.22	19 (3%) 40 26	85, 127, 176, 216	0
All	All	1997/2340 (85%)	0.17	62 (3%) 49 32	63, 116, 179, 250	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	560	ILE	4.8
1	D	115	LEU	4.2
1	C	238	CYS	3.5
1	A	115	LEU	3.5
1	D	532	SER	3.4
1	B	87	ASN	3.3
1	D	571	PHE	3.3
1	C	122	LEU	3.3
1	C	91	ILE	3.3
1	C	98	ARG	3.2
1	D	374	ILE	3.2
1	B	582	VAL	2.9
1	A	123	THR	2.8
1	C	304	ALA	2.8
1	D	354	PHE	2.8
1	B	91	ILE	2.8
1	D	533	LEU	2.7
1	A	103	MET	2.7
1	B	94	LEU	2.6
1	D	444	LEU	2.6
1	D	548	ILE	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	128	LEU	2.5
1	D	342	CYS	2.5
1	D	351	PRO	2.4
1	A	126	THR	2.4
1	C	97	CYS	2.3
1	C	583	GLY	2.3
1	B	561	VAL	2.3
1	A	135	LEU	2.3
1	D	107	LEU	2.3
1	C	107	LEU	2.3
1	D	85	SER	2.3
1	C	561	VAL	2.3
1	B	119	ILE	2.2
1	A	146	VAL	2.2
1	A	170	LYS	2.2
1	B	112	ILE	2.2
1	B	90	VAL	2.2
1	B	304	ALA	2.2
1	D	375	PHE	2.2
1	C	130	LEU	2.2
1	D	105	LEU	2.2
1	B	258	GLY	2.2
1	B	309	LEU	2.2
1	A	238	CYS	2.2
1	A	354	PHE	2.1
1	A	205	THR	2.1
1	B	85	SER	2.1
1	A	119	ILE	2.1
1	A	130	LEU	2.1
1	C	191	LEU	2.1
1	A	125	LEU	2.1
1	B	147	ASN	2.1
1	C	115	LEU	2.1
1	D	401	TRP	2.0
1	D	90	VAL	2.0
1	C	571	PHE	2.0
1	D	120	LYS	2.0
1	C	568	ILE	2.0
1	B	280	ILE	2.0
1	A	143	GLY	2.0
1	D	572	LEU	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 monosaccharides 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.