



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:45 PM UTC

PDB ID : 7EA1 / pdb_00007ea1
Title : Crystal Structure of Spindlin1 bound to SPINDOC Docpep2
Authors : Zhao, F.; Li, H.
Deposited on : 2021-03-05
Resolution : 2.70 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
Xtrriage (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

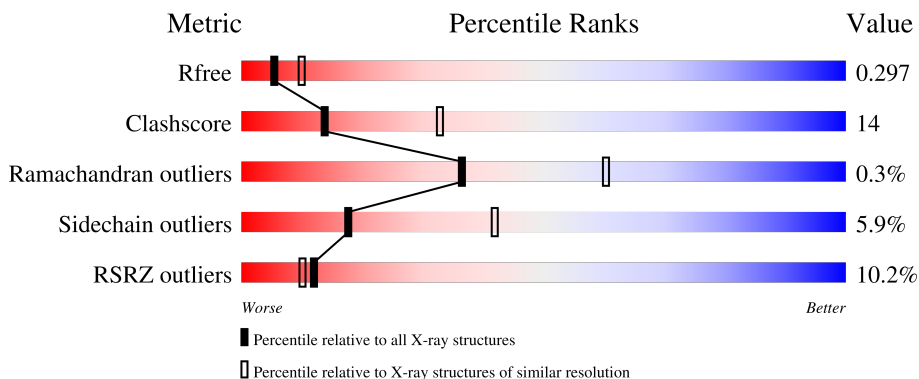
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	215	 4% 72% 20% • 7%
1	C	215	 13% 53% 29% • 14%
2	B	12	 8% 25% 17% 58%
2	D	12	 8% 8% 83%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3188 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 Spindlin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	200	Total 1622	C 1034	N 271	O 308	S 9	0	1	0
1	C	184	Total 1486	C 951	N 246	O 281	S 8	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	48	GLY	-	expression tag	UNP Q9Y657
A	49	SER	-	expression tag	UNP Q9Y657
C	48	GLY	-	expression tag	UNP Q9Y657
C	49	SER	-	expression tag	UNP Q9Y657

- Molecule 2 is a protein called Peptide from Spindlin interactor and repressor of chromatin-binding protein.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	B	5	Total 47	C 29	N 13	O 5	0	0	0
2	D	2	Total 20	C 12	N 6	O 2	0	0	0

- Molecule 3 is water.

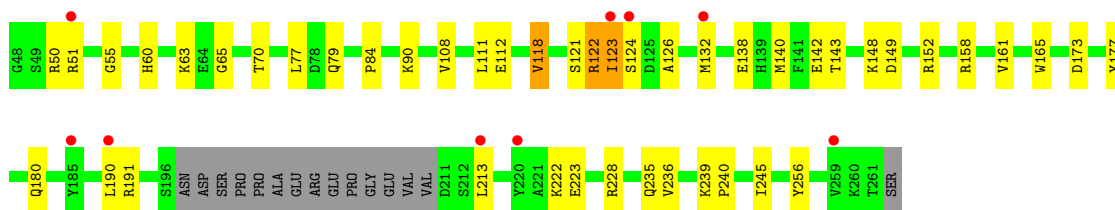
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	10	Total 10	O 10	0	0
3	C	3	Total 3	O 3	0	0

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.

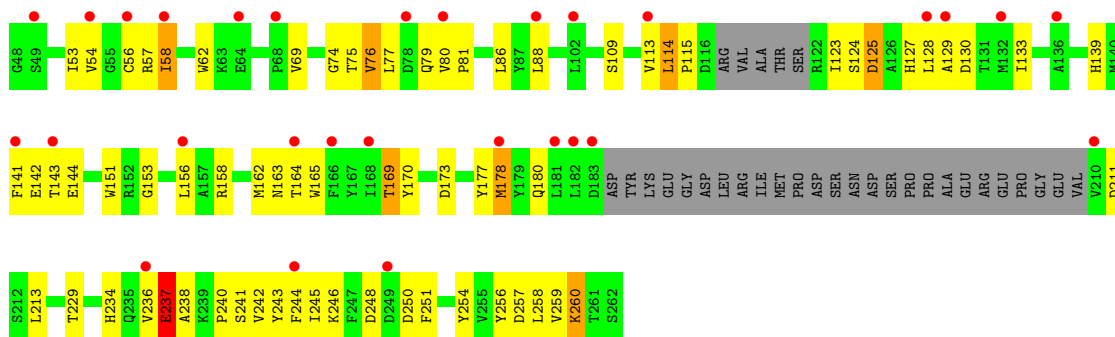
- Molecule 1: Spindlin-1

Chain A: 4% 72% 20% 7%



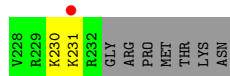
- Molecule 1: Spindlin-1

Chain C: 13% 53% 29% 14%



- Molecule 2: Peptide from Spindlin interactor and repressor of chromatin-binding protein

Chain B: 8% 25% 17% 58%



- Molecule 2: Peptide from Spindlin interactor and repressor of chromatin-binding protein

Chain D: 8% 8% 83%

VAL	
R229	
LYS	
ARG	
GLY	
ARG	
PRO	
MET	
THR	
LYS	
ASN	

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.80Å 136.90Å 42.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.36 – 2.70 40.36 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.3 (40.36-2.70) 99.8 (40.36-2.70)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.14 (at 2.69Å)	Xtrriage
Refinement program	PHENIX 1.14_3247	Depositor
R, R_{free}	0.254 , 0.286 (Not available) , 0.297	Depositor DCC
R_{free} test set	669 reflections (4.72%)	wwPDB-VP
Wilson B-factor (Å ²)	66.1	Xtrriage
Anisotropy	0.718	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 73.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	3188	wwPDB-VP
Average B, all atoms (Å ²)	100.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.01% 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	0.48	3/1663 (0.2%)	0.60	1/2250 (0.0%)
1	C	0.44	1/1520 (0.1%)	0.73	2/2056 (0.1%)
2	B	0.23	0/46	0.80	0/57
2	D	0.19	0/19	0.57	0/22
All	All	0.46	4/3248 (0.1%)	0.67	3/4385 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	239	LYS	C-O	-10.38	1.19	1.23
1	A	121	SER	CA-CB	-5.85	1.45	1.53
1	C	240	PRO	C-O	-5.41	1.17	1.24
1	A	121	SER	C-O	-5.18	1.17	1.24

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	125	ASP	CB-CA-C	6.78	119.82	110.06
1	C	237	GLU	CB-CA-C	-5.45	102.33	110.88
1	A	122	ARG	CB-CA-C	5.10	118.75	109.62

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	1622	0	1577	27	1
1	C	1486	0	1445	66	0
2	B	47	0	60	1	0
2	D	20	0	25	1	0
3	A	10	0	0	0	0
3	C	3	0	0	0	0
All	All	3188	0	3107	91	1

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 14.

All (91) 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:58:ILE:HD11	1:C:76:VAL:CG1	1.99	0.93
1:A:123:ILE:HG12	1:A:158:ARG:HH22	1.35	0.91
1:C:234:HIS:NE2	1:C:236:VAL:HG13	1.86	0.90
1:C:143:THR:HG22	1:C:144:GLU:HG3	1.56	0.87
1:C:234:HIS:NE2	1:C:236:VAL:CG1	2.41	0.83
1:C:77:LEU:HD23	1:C:88:LEU:O	1.80	0.81
1:A:138:GLU:HG3	1:A:191:ARG:HH11	1.45	0.81
1:C:241:SER:HB2	1:C:259:VAL:HG23	1.63	0.79
1:A:122:ARG:HG2	1:A:122:ARG:HH21	1.49	0.77
1:C:58:ILE:HA	1:C:114:LEU:HD22	1.68	0.75
1:C:234:HIS:CD2	1:C:236:VAL:HG13	2.23	0.74
1:A:142:GLU:OE1	1:A:142:GLU:N	2.19	0.72
1:A:173:ASP:OD2	1:A:177:TYR:OH	2.10	0.70
1:C:77:LEU:HD23	1:C:77:LEU:H	1.58	0.69
1:C:58:ILE:HD11	1:C:76:VAL:HG11	1.75	0.67
1:C:141:PHE:HZ	2:D:229:ARG:HH11	1.42	0.67
1:A:122:ARG:HG2	1:A:122:ARG:NH2	2.10	0.66
2:B:230:LYS:HG3	1:C:211:ASP:HB2	1.76	0.66
1:C:58:ILE:HD11	1:C:76:VAL:HG13	1.78	0.65
1:C:75:THR:HG21	1:C:162:MET:HE2	1.78	0.65
1:C:54:VAL:HA	1:C:76:VAL:HG23	1.79	0.65
1:C:58:ILE:CD1	1:C:76:VAL:CG1	2.75	0.64
1:C:54:VAL:HA	1:C:76:VAL:CG2	2.28	0.63
1:C:153:GLY:HA3	1:C:170:TYR:HA	1.82	0.62
1:C:77:LEU:HD23	1:C:77:LEU:N	2.16	0.60
1:C:241:SER:HB2	1:C:259:VAL:CG2	2.31	0.60
1:C:58:ILE:N	1:C:74:GLY:O	2.25	0.59
1:C:123:ILE:HG22	1:C:123:ILE:O	2.01	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:234:HIS:NE2	1:C:236:VAL:HG12	2.19	0.58
1:C:56:CYS:HB2	1:C:114:LEU:HB2	1.85	0.57
1:C:142:GLU:HG2	1:C:142:GLU:O	2.04	0.57
1:C:250:ASP:OD1	1:C:251:PHE:N	2.38	0.56
1:C:58:ILE:HD13	1:C:114:LEU:HD21	1.89	0.54
1:C:77:LEU:HD23	1:C:88:LEU:C	2.33	0.54
1:C:123:ILE:HG22	1:C:125:ASP:OD1	2.07	0.53
1:C:77:LEU:N	1:C:77:LEU:CD2	2.71	0.53
1:C:237:GLU:N	1:C:237:GLU:OE2	2.42	0.53
1:A:77:LEU:HD11	1:A:90:LYS:HB2	1.91	0.51
1:C:56:CYS:HB2	1:C:114:LEU:HD12	1.92	0.50
1:C:133:ILE:HG22	1:C:238:ALA:HB2	1.91	0.50
1:A:122:ARG:NH2	1:A:122:ARG:CG	2.72	0.50
1:C:77:LEU:CD2	1:C:88:LEU:O	2.57	0.50
1:A:213:LEU:H	1:A:213:LEU:HD23	1.77	0.50
1:C:139:HIS:O	1:C:151:TRP:N	2.39	0.49
1:C:162:MET:HG3	1:C:178:MET:HE1	1.95	0.49
1:C:77:LEU:CD2	1:C:88:LEU:C	2.87	0.48
1:C:162:MET:HB3	1:C:165:TRP:CD1	2.48	0.48
1:C:242:VAL:HA	1:C:256:TYR:O	2.12	0.48
1:A:55:GLY:HA2	1:A:161:VAL:HG11	1.95	0.48
1:C:173:ASP:OD2	1:C:177:TYR:OH	2.31	0.48
1:C:250:ASP:OD2	1:C:254:TYR:OH	2.32	0.47
1:A:60:HIS:CG	1:A:108:VAL:HG13	2.49	0.47
1:C:53:ILE:HG13	1:C:76:VAL:HG21	1.96	0.47
1:C:123:ILE:O	1:C:125:ASP:OD1	2.32	0.47
1:C:127:HIS:O	1:C:158:ARG:NH1	2.49	0.46
1:C:129:ALA:HB2	1:C:158:ARG:NH2	2.31	0.46
1:C:243:TYR:CD1	1:C:258:LEU:HD12	2.50	0.46
1:C:62:TRP:H	1:C:69:VAL:HG13	1.81	0.46
1:C:129:ALA:HB2	1:C:158:ARG:HH22	1.80	0.45
1:A:50:ARG:O	1:A:51:ARG:HG3	2.16	0.45
1:A:132:MET:HE2	1:A:190:LEU:HD23	1.98	0.45
1:A:70:THR:HG21	1:C:213:LEU:HD13	1.98	0.45
1:C:130:ASP:N	1:C:130:ASP:OD1	2.48	0.45
1:C:169:THR:OG1	1:C:170:TYR:N	2.50	0.45
1:C:163:ASN:OD1	1:C:163:ASN:N	2.49	0.45
1:A:143:THR:HG23	1:A:149:ASP:HB2	1.99	0.44
1:A:245:ILE:HD12	1:A:256:TYR:HD2	1.82	0.44
1:C:57:ARG:NH1	1:C:75:THR:OG1	2.51	0.44
1:C:245:ILE:HD13	1:C:256:TYR:HD1	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:123:ILE:HG21	1:A:126:ALA:HA	2.00	0.43
1:C:229:THR:OG1	1:C:248:ASP:OD2	2.34	0.43
1:A:138:GLU:OE1	1:A:152:ARG:HD3	2.18	0.43
1:A:123:ILE:H	1:A:123:ILE:HG13	1.41	0.43
1:A:79:GLN:NE2	1:A:84:PRO:O	2.34	0.43
1:A:235:GLN:NE2	1:A:240:PRO:O	2.46	0.43
1:C:80:VAL:HG21	1:C:86:LEU:HD22	1.99	0.42
1:C:133:ILE:HG23	1:C:156:LEU:O	2.19	0.42
1:A:60:HIS:HB3	1:A:111:LEU:HA	2.02	0.42
1:C:242:VAL:HG12	1:C:257:ASP:HB3	2.01	0.42
1:C:58:ILE:CD1	1:C:76:VAL:HG13	2.45	0.42
1:C:164:THR:O	1:C:180:GLN:HB2	2.20	0.41
1:A:65:GLY:HA3	1:C:260:LYS:HB3	2.02	0.41
1:C:246:LYS:C	1:C:246:LYS:HD3	2.46	0.41
1:C:58:ILE:CG1	1:C:76:VAL:HG13	2.51	0.41
1:C:113:VAL:C	1:C:114:LEU:HD23	2.46	0.41
1:A:140:MET:SD	1:A:148:LYS:HB3	2.60	0.41
1:A:165:TRP:CE2	1:A:180:GLN:HG3	2.56	0.40
1:C:156:LEU:HD13	1:C:244:PHE:CE1	2.55	0.40
1:A:55:GLY:O	1:A:118:VAL:HG13	2.21	0.40
1:A:123:ILE:HG21	1:A:158:ARG:HH12	1.86	0.40
1:C:79:GLN:HG3	1:C:81:PRO:HD3	2.04	0.40

All (1) 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)
1:A:112:GLU:OE2	1:A:122:ARG:NH1[2_554]	1.99	0.21

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	129/215 (60%)	123 (95%)	6 (5%)	0	100	100
1	C	178/215 (83%)	161 (90%)	16 (9%)	1 (1%)	21	44
2	B	3/12 (25%)	1 (33%)	2 (67%)	0	100	100
All	All	310/442 (70%)	285 (92%)	24 (8%)	1 (0%)	36	60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	115	PRO

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	177/189 (94%)	169 (96%)	8 (4%)	24	52
1	C	139/189 (74%)	129 (93%)	10 (7%)	13	32
2	B	5/11 (46%)	4 (80%)	1 (20%)	1	4
2	D	2/11 (18%)	2 (100%)	0	100	100
All	All	323/400 (81%)	304 (94%)	19 (6%)	18	42

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	LYS
1	A	118	VAL
1	A	123	ILE
1	A	124	SER
1	A	222	LYS
1	A	223	GLU
1	A	228	ARG
1	A	236	VAL
2	B	231	LYS
1	C	58	ILE
1	C	76	VAL

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Mol	Chain	Res	Type
1	C	109	SER
1	C	114	LEU
1	C	124	SER
1	C	128	LEU
1	C	169	THR
1	C	178	MET
1	C	237	GLU
1	C	260	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	GLN
1	A	103	ASN
1	A	163	ASN

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	200/215 (93%)	0.48	9 (4%) 38 34	31, 66, 143, 217	1 (0%)
1	C	184/215 (85%)	1.02	29 (15%) 5 4	58, 122, 187, 225	0
2	B	5/12 (41%)	1.85	1 (20%) 3 2	136, 157, 170, 177	0
2	D	2/12 (16%)	1.46	1 (50%) 0 0	134, 134, 134, 154	0
All	All	391/454 (86%)	0.76	40 (10%) 12 10	31, 90, 179, 225	1 (0%)

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	123	ILE	4.9
1	A	185	TYR	4.6
1	C	210	VAL	4.5
1	C	68	PRO	4.0
1	A	220	TYR	3.5
1	C	102	LEU	3.2
1	A	259	VAL	3.2
1	C	49	SER	3.1
1	C	182	LEU	3.0
1	C	129	ALA	3.0
1	C	136	ALA	2.9
1	C	132	MET	2.8
1	C	156	LEU	2.8
1	C	113	VAL	2.6
1	C	64	GLU	2.5
2	B	231	LYS	2.5
1	A	51	ARG	2.5
1	A	124	SER	2.5
1	C	80	VAL	2.5
1	A	213	LEU	2.4
1	C	88	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	78	ASP	2.4
1	C	236	VAL	2.4
1	C	166	PHE	2.4
1	A	132	MET	2.3
1	C	183	ASP	2.3
1	C	249	ASP	2.3
1	C	58	ILE	2.2
1	C	168	ILE	2.2
1	C	164	THR	2.2
1	A	190	LEU	2.2
1	C	178	MET	2.2
1	C	244	PHE	2.2
1	C	56	CYS	2.1
1	C	128	LEU	2.1
1	C	143	THR	2.1
1	C	141	PHE	2.1
1	C	181	LEU	2.0
1	C	54	VAL	2.0
2	D	229	ARG	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.