Full wwPDB X-ray Structure Validation Report (i)

Jun 12, 2024 – 02:56 PM EDT

PDB ID	:	3W6J
Title	:	Crystal structure of ScpAB core complex
Authors	:	Kamada, K.; Hirano, T.
Deposited on	:	2013-02-15
Resolution	:	2.60 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 (i) 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 (i)) were used in the production of this report:

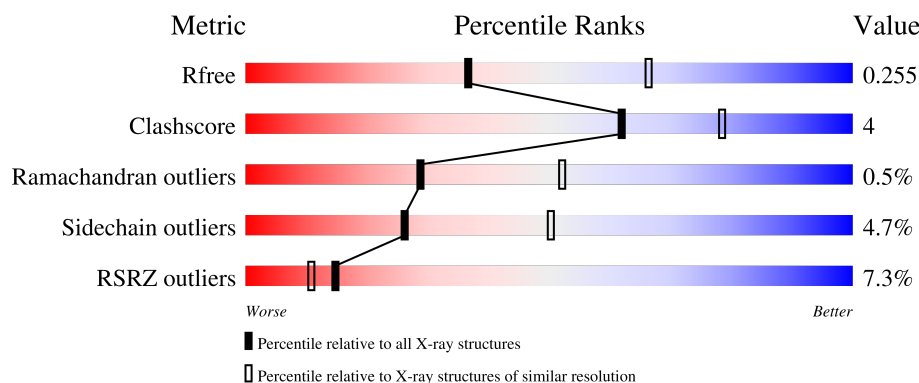|		
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 2.60 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	Percentile Ranks	Value
Rfree		0.255
Clashscore		4
Ramachandran outliers		0.5%
Sidechain outliers		4.7%
RSRZ outliers		7.3%

Worse — *Better*

▮ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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 <=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	174	% · 74% 12% 14%
1	D	174	6% · 64% 13% · 21%
2	B	184	14% · 70% 20% · 9%
2	C	184	3% · 85% 6% 9%
2	E	184	9% · 70% 18% · 11%

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Mol	Chain	Length	Quality of chain
2	F	184	

2 Entry composition ⓘ

There are 3 unique types of molecules in this entry. The entry contains 7542 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 ScpA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	150	Total	C	N	O	S	0	0	0
			1230	796	202	225	7			
1	D	137	Total	C	N	O	S	0	0	0
			1117	722	186	203	6			

- Molecule 2 is a protein called ScpB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	168	Total	C	N	O	S	0	0	0
			1294	831	216	245	2			
2	C	167	Total	C	N	O	S	0	0	0
			1279	821	214	242	2			
2	E	164	Total	C	N	O	S	0	0	0
			1272	818	212	240	2			
2	F	170	Total	C	N	O	S	0	0	0
			1307	840	218	247	2			

- Molecule 3 is water.

| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	9	Total	O	0	0
			9	9		
3	B	3	Total	O	0	0
			3	3		
3	C	16	Total	O	0	0
			16	16		
3	D	5	Total	O	0	0
			5	5		
3	F	10	Total	O	0	0
			10	10		

3 Residue-property plots ⓘ

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.
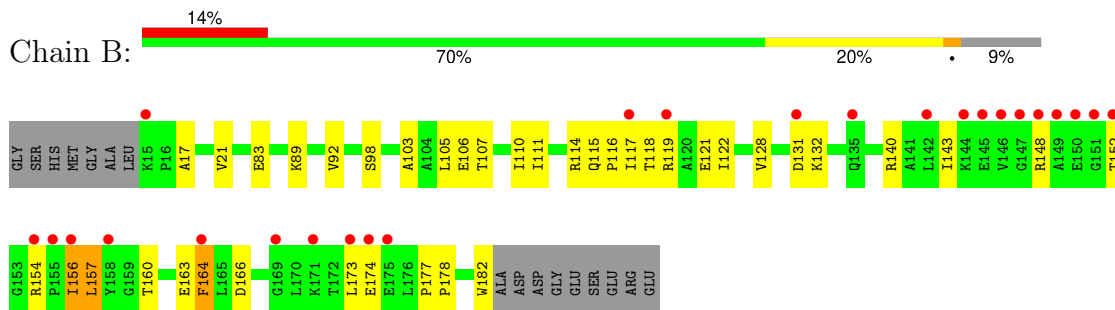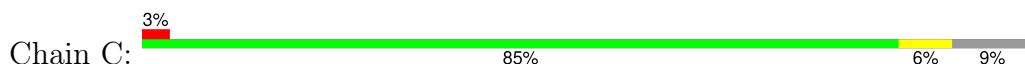
- Molecule 1: ScpA
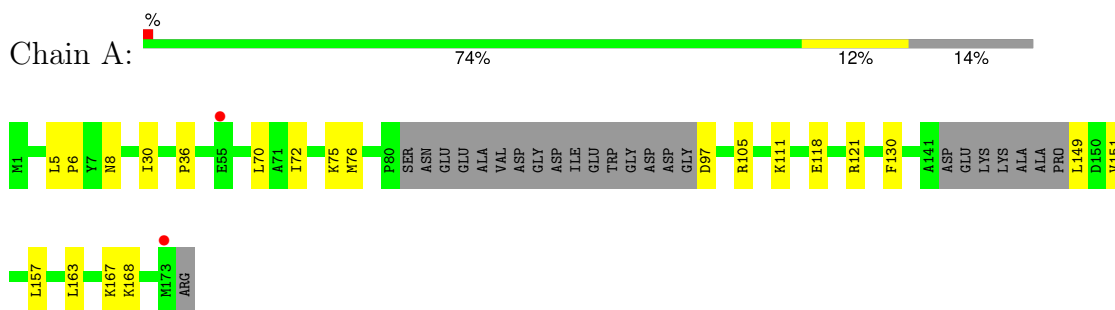
- Molecule 1: ScpA
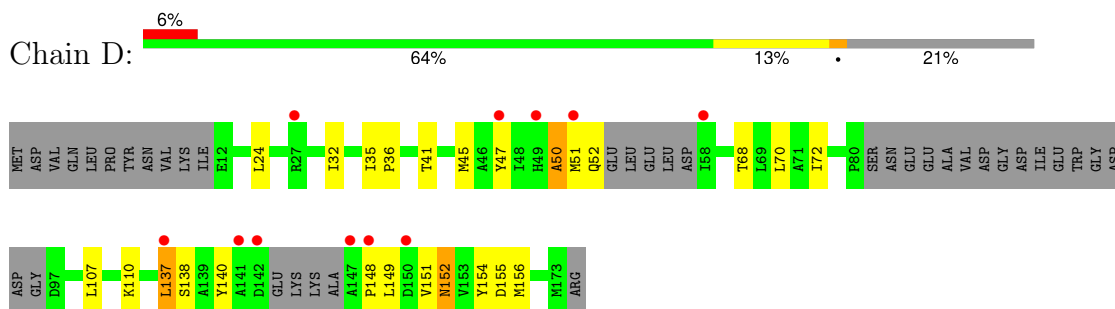
- Molecule 2: ScpB

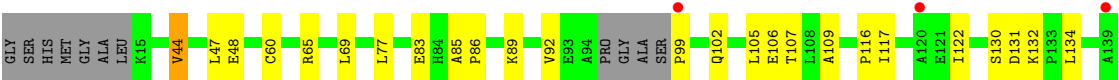- Molecule 2: ScpB

• Molecule 2: ScpB

Chain E:

• Molecule 2: ScpB

Chain F:
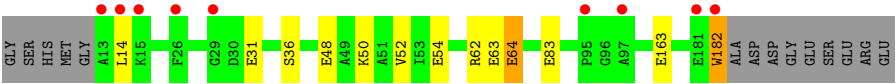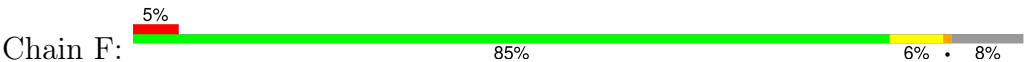
4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	90.86Å 127.42Å 97.22Å 90.00° 113.84° 90.00°	Depositor
Resolution (Å)	30.70 – 2.60 30.70 – 2.59	Depositor EDS
% Data completeness (in resolution range)	98.7 (30.70-2.60) 98.7 (30.70-2.59)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$< I/\sigma(I) >$ [1]	2.11 (at 2.61Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.225 , 0.264 0.216 , 0.255	Depositor DCC
R_{free} test set	3145 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å²)	53.9	Xtriage
Anisotropy	0.502	Xtriage
Bulk solvent k_{sol}(e/Å³), B_{sol}(Å²)	0.34 , 44.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	7542	wwPDB-VP
Average B, all atoms (Å²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.63% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< \|L\| >$, $< L^2 >$ 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	0.45	0/1250	0.57	0/1684
1	D	0.41	0/1135	0.57	0/1527
2	B	0.42	0/1315	0.53	0/1781
2	C	0.44	0/1298	0.63	0/1757
2	E	0.40	0/1291	0.54	0/1745
2	F	0.46	0/1328	0.58	0/1799
All	All	0.43	0/7617	0.57	0/10293

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 ⓘ

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	1230	0	1281	14	0
1	D	1117	0	1159	16	0
2	B	1294	0	1347	20	0
2	C	1279	0	1342	8	0
2	E	1272	0	1327	17	0
2	F	1307	0	1363	6	0
3	A	9	0	0	1	0
3	B	3	0	0	0	0
3	C	16	0	0	0	0
3	D	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	F	10	0	0	0	0
All	All	7542	0	7819	68	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 (68) 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)
2:B:111:ILE:HG21	2:B:143:ILE:HD11	1.72	0.71
1:A:8:ASN:HB3	2:E:132:LYS:HD3	1.79	0.65
2:F:62:ARG:NH1	2:F:64:GLU:OE2	2.34	0.61
2:C:58:GLN:OE1	2:C:61:ARG:NH1	2.34	0.60
2:B:118:THR:HB	2:B:157:LEU:HD22	1.86	0.58
1:D:152:ASN:ND2	1:D:155:ASP:OD2	2.40	0.55
1:D:149:LEU:HB2	1:D:151:VAL:HG23	1.89	0.55
2:F:14:LEU:HD22	2:F:52:VAL:HG22	1.89	0.54
1:A:130:PHE:CE2	1:A:167:LYS:HD2	2.42	0.54
1:A:168:LYS:HB2	2:C:94:ALA:HB3	1.89	0.53
1:D:149:LEU:HD21	2:E:99:PRO:HG2	1.90	0.53
2:E:148:ARG:NH1	2:E:154:ARG:O	2.34	0.52
2:B:17:ALA:O	2:B:21:VAL:HG23	2.09	0.52
2:B:116:PRO:HB2	2:B:157:LEU:HD12	1.91	0.52
1:D:50:ALA:O	1:D:52:GLN:N	2.42	0.52
1:A:151:VAL:HG11	2:B:105:LEU:HD21	1.92	0.51
2:B:111:ILE:O	2:B:115:GLN:N	2.43	0.51
2:E:171:LYS:HB2	2:E:175:GLU:OE2	2.11	0.51
2:E:130:SER:O	2:E:134:LEU:HB2	2.11	0.51
1:A:105:ARG:NH1	3:A:209:HOH:O	2.44	0.50
1:D:41:THR:O	1:D:45:MET:HG2	2.11	0.50
2:B:160:THR:HB	2:B:164:PHE:CD2	2.47	0.50
2:B:103:ALA:O	2:B:107:THR:OG1	2.24	0.50
1:A:105:ARG:NE	2:C:124:GLU:OE2	2.45	0.49
2:E:102:GLN:O	2:E:106:GLU:HG3	2.12	0.49
2:B:148:ARG:NH1	2:B:154:ARG:O	2.42	0.49
2:B:106:GLU:O	2:B:110:ILE:HG13	2.13	0.48
2:F:48:GLU:O	2:F:52:VAL:HG23	2.13	0.48
2:B:111:ILE:HD13	2:B:143:ILE:HG12	1.96	0.48
2:B:92:VAL:HA	2:B:98:SER:HB3	1.95	0.48
1:D:156:MET:HE3	2:E:164:PHE:HE1	1.78	0.48
1:A:118:GLU:OE1	1:A:121:ARG:NH1	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:156:MET:HG2	2:E:109:ALA:HB2	1.97	0.47
1:D:152:ASN:ND2	1:D:154:TYR:HB3	2.30	0.46
1:A:72:ILE:O	1:A:76:MET:HG3	2.15	0.46
2:F:50:LYS:HE2	2:F:54:GLU:HG3	1.98	0.45
1:A:5:LEU:HA	1:A:6:PRO:HD3	1.69	0.45
1:A:157:LEU:HD11	2:B:177:PRO:HG2	1.99	0.45
1:A:163:LEU:HD13	2:B:106:GLU:OE1	2.16	0.45
2:E:44:VAL:HG13	2:E:48:GLU:HB3	1.99	0.45
1:D:24:LEU:HD21	1:D:47:TYR:CE1	2.52	0.45
2:B:110:ILE:O	2:B:114:ARG:HB2	2.16	0.44
1:A:111:LYS:HD3	2:C:179:LEU:HG	2.00	0.44
2:E:116:PRO:HB2	2:E:157:LEU:HD13	1.98	0.44
1:D:110:LYS:HD3	2:F:182:TRP:CZ2	2.52	0.43
2:E:86:PRO:HA	2:E:89:LYS:HE2	1.99	0.43
2:B:89:LYS:HE3	2:B:89:LYS:HB2	1.82	0.43
2:E:102:GLN:HB2	2:F:31:GLU:OE2	2.19	0.43
2:E:117:ILE:HG12	2:E:122:ILE:HG13	2.01	0.43
2:E:60:CYS:HA	2:E:65:ARG:HG3	2.00	0.43
2:E:107:THR:HG23	2:E:122:ILE:HG21	2.01	0.43
1:D:70:LEU:HD23	1:D:70:LEU:HA	1.82	0.42
1:A:70:LEU:HD23	1:A:70:LEU:HA	1.85	0.42
2:B:118:THR:O	2:B:121:GLU:N	2.52	0.42
1:A:30:ILE:HG23	1:A:36:PRO:HD2	2.00	0.42
1:D:68:THR:O	1:D:72:ILE:HG13	2.19	0.42
2:E:85:ALA:N	2:E:86:PRO:HD2	2.34	0.42
1:D:156:MET:HG3	2:E:105:LEU:HD22	2.01	0.41
1:D:35:ILE:HA	1:D:36:PRO:HD2	1.91	0.41
2:C:145:GLU:HG2	2:C:156:ILE:HG21	2.03	0.41
2:B:177:PRO:HA	2:B:178:PRO:HD2	1.90	0.41
2:C:148:ARG:NH2	2:C:156:ILE:HD11	2.35	0.41
1:D:32:ILE:HD13	1:D:107:LEU:HD23	2.03	0.41
1:D:137:LEU:HB3	1:D:140:TYR:HD1	1.86	0.41
2:B:117:ILE:HD13	2:B:122:ILE:HG12	2.02	0.41
2:C:179:LEU:HD12	2:C:179:LEU:HA	1.88	0.41
2:C:36:SER:HB3	2:C:46:GLU:OE2	2.21	0.41
2:B:156:ILE:H	2:B:156:ILE:HD12	1.86	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	144/174 (83%)	138 (96%)	6 (4%)	0	100	100
1	D	129/174 (74%)	117 (91%)	9 (7%)	3 (2%)	6	11
2	B	166/184 (90%)	151 (91%)	13 (8%)	2 (1%)	13	27
2	C	165/184 (90%)	165 (100%)	0	0	100	100
2	E	160/184 (87%)	147 (92%)	13 (8%)	0	100	100
2	F	168/184 (91%)	166 (99%)	2 (1%)	0	100	100
All	All	932/1084 (86%)	884 (95%)	43 (5%)	5 (0%)	29	52

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	152	THR
1	D	51	MET
2	B	119	ARG
1	D	50	ALA
1	D	148	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	134/152 (88%)	131 (98%)	3 (2%)	52	76
1	D	120/152 (79%)	117 (98%)	3 (2%)	47	73
2	B	135/146 (92%)	122 (90%)	13 (10%)	8	16

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	C	134/146 (92%)	133 (99%)	1 (1%)	84	94
2	E	133/146 (91%)	122 (92%)	11 (8%)	11	22
2	F	136/146 (93%)	130 (96%)	6 (4%)	28	53
All	All	792/888 (89%)	755 (95%)	37 (5%)	26	50

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	75	LYS
1	A	97	ASP
1	A	149	LEU
2	B	83	GLU
2	B	128	VAL
2	B	131	ASP
2	B	132	LYS
2	B	140	ARG
2	B	156	ILE
2	B	157	LEU
2	B	163	GLU
2	B	164	PHE
2	B	166	ASP
2	B	173	LEU
2	B	174	GLU
2	B	182	TRP
2	C	55	GLU
1	D	137	LEU
1	D	138	SER
1	D	152	ASN
2	E	44	VAL
2	E	47	LEU
2	E	69	LEU
2	E	77	LEU
2	E	83	GLU
2	E	92	VAL
2	E	131	ASP
2	E	143	ILE
2	E	152	THR
2	E	163	GLU
2	E	168	PHE
2	F	36	SER
2	F	63	GLU

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Mol	Chain	Res	Type
2	F	64	GLU
2	F	83	GLU
2	F	163	GLU
2	F	182	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	152	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

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	150/174 (86%)	-0.18	2 (1%)	77	73	38, 56, 92, 114	0
1	D	137/174 (78%)	0.27	11 (8%)	12	9	42, 68, 110, 126	0
2	B	168/184 (91%)	0.57	25 (14%)	2	1	40, 87, 125, 136	0
2	C	167/184 (90%)	-0.07	6 (3%)	42	35	34, 48, 74, 99	0
2	E	164/184 (89%)	0.40	17 (10%)	6	4	42, 82, 116, 129	0
2	F	170/184 (92%)	-0.01	9 (5%)	26	20	38, 51, 83, 112	0
All	All	956/1084 (88%)	0.16	70 (7%)	15	11	34, 61, 113, 136	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	167	TYR	7.1
2	B	150	GLU	5.6
1	D	51	MET	5.3
2	E	164	PHE	5.2
1	D	148	PRO	4.8
2	B	154	ARG	4.7
1	D	147	ALA	4.4
2	B	15	LYS	4.2
2	B	152	THR	4.1
2	B	173	LEU	4.1
2	E	149	ALA	4.1
2	E	154	ARG	4.0
2	E	150	GLU	3.9
2	B	147	GLY	3.9
2	E	151	GLY	3.8
2	B	117	ILE	3.7
2	B	146	VAL	3.7
2	B	156	ILE	3.7
2	F	13	ALA	3.6

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Mol	Chain	Res	Type	RSRZ
2	B	144	LYS	3.4
2	B	158	TYR	3.4
2	E	143	ILE	3.3
2	E	152	THR	3.2
1	D	47	TYR	3.2
2	F	181	GLU	3.2
2	C	15	LYS	3.2
2	E	153	GLY	3.2
2	E	171	LYS	3.2
1	D	49	HIS	3.1
1	D	150	ASP	3.1
1	D	58	ILE	3.0
2	F	15	LYS	2.9
1	A	55	GLU	2.9
2	B	142	LEU	2.9
2	F	182	TRP	2.8
2	C	150	GLU	2.8
2	B	175	GLU	2.8
2	E	99	PRO	2.7
2	F	14	LEU	2.7
1	D	137	LEU	2.7
2	B	164	PHE	2.6
2	C	152	THR	2.6
2	B	155	PRO	2.6
2	B	119	ARG	2.6
2	B	145	GLU	2.5
2	B	131	ASP	2.5
2	B	135	GLN	2.5
2	B	174	GLU	2.5
2	E	168	PHE	2.4
2	B	148	ARG	2.4
2	B	151	GLY	2.4
2	E	120	ALA	2.4
2	E	175	GLU	2.4
2	F	95	PRO	2.3
2	E	163	GLU	2.3
2	E	172	THR	2.3
1	D	27	ARG	2.2
2	E	139	ALA	2.2
2	F	29	GLY	2.2
1	D	142	ASP	2.1
2	B	149	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
2	F	97	ALA	2.1
1	A	173	MET	2.1
2	C	29	GLY	2.1
1	D	141	ALA	2.1
2	C	14	LEU	2.0
2	B	171	LYS	2.0
2	F	26	PHE	2.0
2	C	28	ALA	2.0
2	B	169	GLY	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 monosaccharides in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.