



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2017 – 05:36 PM EDT

PDB ID : 5XEI
Title : Crystal structure of the Smc head domain with a coiled coil and joint derived from *Pyrococcus yayanosii*
Authors : Lee, H.; Noh, H.; Oh, B.-H.
Deposited on : unknown
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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029824
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20029824

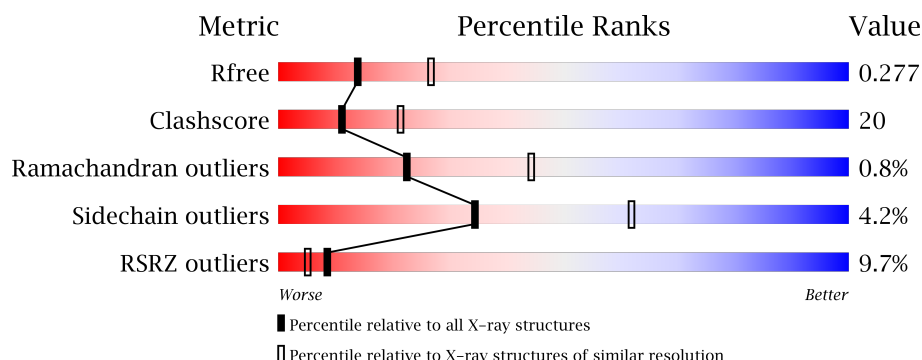
1 Overall quality at a glance

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	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	2542 (2.60-2.60)
Clashscore	112137	2895 (2.60-2.60)
Ramachandran outliers	110173	2848 (2.60-2.60)
Sidechain outliers	110143	2848 (2.60-2.60)
RSRZ outliers	101464	2550 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	540	<div> <div>9%</div> <div>64%</div> <div>30%</div> <div>••</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4099 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 Chromosome partition protein Smc.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	517	Total	C	N	O	S	0	0	0
			4086	2590	707	776	13			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	903	SER	-	linker	UNP F8AFS8
A	904	GLY	-	linker	UNP F8AFS8
A	905	GLY	-	linker	UNP F8AFS8
A	906	SER	-	linker	UNP F8AFS8
A	1178	VAL	-	expression tag	UNP F8AFS8
A	1179	ASP	-	expression tag	UNP F8AFS8
A	1180	ALA	-	expression tag	UNP F8AFS8
A	1181	GLY	-	expression tag	UNP F8AFS8
A	1182	SER	-	expression tag	UNP F8AFS8
A	1183	SER	-	expression tag	UNP F8AFS8
A	1184	SER	-	expression tag	UNP F8AFS8
A	1185	ARG	-	expression tag	UNP F8AFS8
A	1186	PRO	-	expression tag	UNP F8AFS8
A	1187	GLY	-	expression tag	UNP F8AFS8
A	1188	LEU	-	expression tag	UNP F8AFS8

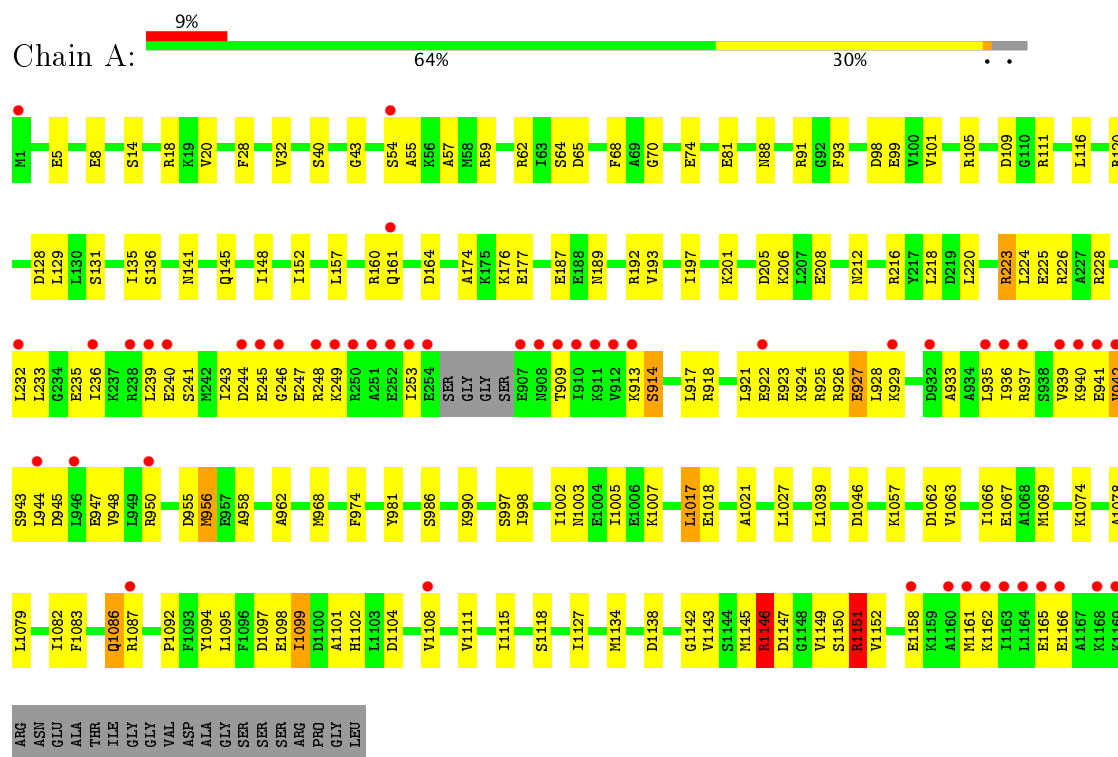
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	13	Total	O	0	0
			13	13		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: Chromosome partition protein Smc



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	120.32Å 50.84Å 122.06Å 90.00° 118.09° 90.00°	Depositor
Resolution (Å)	39.62 – 2.60 39.63 – 2.60	Depositor EDS
% Data completeness (in resolution range)	92.6 (39.62-2.60) 92.6 (39.63-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.73 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.207 , 0.278 0.205 , 0.277	Depositor DCC
R_{free} test set	1902 reflections (10.08%)	DCC
Wilson B-factor (Å ²)	50.9	Xtriage
Anisotropy	0.189	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 56.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.028 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4099	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.96% 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	0.53	0/4135	0.69	3/5537 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	1151	ARG	NE-CZ-NH1	-11.44	114.58	120.30
1	A	1151	ARG	NE-CZ-NH2	6.63	123.61	120.30
1	A	1146	ARG	CG-CD-NE	6.48	125.40	111.80

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	4086	0	4206	166	1
2	A	13	0	0	0	0
All	All	4099	0	4206	166	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 20.

All (166) 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:1146:ARG:HD3	1:A:1151:ARG:CZ	1.35	1.52
1:A:1146:ARG:O	1:A:1151:ARG:NH2	1.81	1.12
1:A:1146:ARG:CD	1:A:1151:ARG:CZ	2.30	1.08
1:A:1151:ARG:N	1:A:1151:ARG:HD2	1.70	1.03
1:A:1146:ARG:HD3	1:A:1151:ARG:NE	1.74	1.03
1:A:1146:ARG:H	1:A:1151:ARG:NH2	1.59	1.01
1:A:1146:ARG:HD3	1:A:1151:ARG:NH2	1.74	1.00
1:A:1146:ARG:N	1:A:1151:ARG:HH12	1.66	0.94
1:A:1146:ARG:C	1:A:1151:ARG:HH22	1.71	0.93
1:A:1146:ARG:CA	1:A:1151:ARG:HH22	1.82	0.91
1:A:1146:ARG:N	1:A:1151:ARG:HH22	1.67	0.90
1:A:1146:ARG:H	1:A:1151:ARG:CZ	1.86	0.89
1:A:1146:ARG:H	1:A:1151:ARG:HH22	1.17	0.88
1:A:1098:GLU:HG2	1:A:1099:ILE:H	1.41	0.85
1:A:161:GLN:OE1	1:A:161:GLN:N	2.09	0.84
1:A:923:GLU:OE2	1:A:926:ARG:NH1	2.14	0.81
1:A:1146:ARG:N	1:A:1151:ARG:NH1	2.32	0.78
1:A:236:ILE:HG23	1:A:921:LEU:HD21	1.63	0.77
1:A:1145:MET:HG3	1:A:1149:VAL:O	1.85	0.76
1:A:1146:ARG:H	1:A:1151:ARG:NH1	1.85	0.75
1:A:160:ARG:NH1	1:A:164:ASP:OD2	2.19	0.75
1:A:40:SER:HB2	1:A:59:ARG:HH12	1.53	0.73
1:A:1098:GLU:HG2	1:A:1099:ILE:N	2.03	0.72
1:A:136:SER:HA	1:A:141:ASN:OD1	1.89	0.72
1:A:1146:ARG:HH11	1:A:1151:ARG:HH21	1.38	0.70
1:A:1146:ARG:CD	1:A:1151:ARG:NH2	2.52	0.70
1:A:245:GLU:HA	1:A:248:ARG:HB3	1.73	0.70
1:A:152:ILE:HG13	1:A:1074:LYS:HG3	1.72	0.70
1:A:1143:VAL:HG22	1:A:1152:VAL:HG12	1.74	0.69
1:A:174:ALA:O	1:A:177:GLU:HG3	1.95	0.66
1:A:1066:ILE:HD12	1:A:1067:GLU:N	2.12	0.65
1:A:148:ILE:O	1:A:152:ILE:HG12	1.97	0.65
1:A:128:ASP:OD2	1:A:129:LEU:N	2.30	0.65
1:A:201:LYS:NZ	1:A:205:ASP:OD1	2.30	0.64
1:A:145:GLN:HG2	1:A:1098:GLU:OE1	1.97	0.64
1:A:1146:ARG:CD	1:A:1151:ARG:NE	2.54	0.64
1:A:18:ARG:NH2	1:A:74:GLU:OE1	2.31	0.63
1:A:223:ARG:HH21	1:A:226:ARG:HB2	1.64	0.63
1:A:1146:ARG:NH1	1:A:1149:VAL:HB	2.14	0.63
1:A:247:GLU:HG2	1:A:914:SER:HB2	1.81	0.62
1:A:947:GLU:OE1	1:A:950:ARG:NE	2.32	0.61
1:A:1098:GLU:CG	1:A:1099:ILE:H	2.03	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1146:ARG:HG2	1:A:1146:ARG:O	2.00	0.59
1:A:212:ASN:O	1:A:216:ARG:HG3	2.02	0.59
1:A:223:ARG:HA	1:A:223:ARG:HE	1.67	0.59
1:A:208:GLU:CG	1:A:974:PHE:HE2	2.15	0.58
1:A:1162:LYS:O	1:A:1166:GLU:OE1	2.21	0.58
1:A:1079:LEU:HD21	1:A:1099:ILE:HD12	1.86	0.58
1:A:1146:ARG:CG	1:A:1151:ARG:NH2	2.67	0.57
1:A:1146:ARG:CA	1:A:1151:ARG:NH2	2.61	0.57
1:A:223:ARG:NH2	1:A:226:ARG:HD2	2.19	0.57
1:A:218:LEU:HD13	1:A:968:MET:HE3	1.86	0.56
1:A:40:SER:CB	1:A:59:ARG:HH12	2.17	0.56
1:A:223:ARG:HH21	1:A:226:ARG:HD2	1.70	0.56
1:A:922:GLU:HA	1:A:925:ARG:CZ	2.36	0.56
1:A:32:VAL:HG11	1:A:1134:MET:HG3	1.88	0.55
1:A:943:SER:OG	1:A:944:LEU:N	2.36	0.55
1:A:218:LEU:HD13	1:A:968:MET:CE	2.36	0.55
1:A:176:LYS:HB2	1:A:1005:ILE:HG21	1.88	0.55
1:A:1078:ALA:O	1:A:1082:ILE:HG13	2.06	0.55
1:A:218:LEU:CD1	1:A:968:MET:CE	2.85	0.55
1:A:917:LEU:O	1:A:921:LEU:HB2	2.07	0.55
1:A:91:ARG:NH2	1:A:98:ASP:OD1	2.35	0.55
1:A:1146:ARG:NH1	1:A:1151:ARG:HE	2.04	0.54
1:A:1086:GLN:HB2	1:A:1094:TYR:OH	2.08	0.54
1:A:239:LEU:HD13	1:A:921:LEU:HD23	1.90	0.53
1:A:239:LEU:HB3	1:A:921:LEU:HD23	1.89	0.53
1:A:1097:ASP:HA	1:A:1127:ILE:HB	1.91	0.52
1:A:933:ALA:HA	1:A:936:ILE:HG22	1.92	0.52
1:A:160:ARG:HG2	1:A:160:ARG:HH11	1.73	0.52
1:A:225:GLU:OE2	1:A:228:ARG:HD2	2.10	0.52
1:A:232:LEU:HA	1:A:235:GLU:HB2	1.91	0.52
1:A:236:ILE:HD11	1:A:924:LYS:O	2.10	0.52
1:A:201:LYS:HB2	1:A:981:TYR:CE1	2.46	0.51
1:A:208:GLU:HG3	1:A:974:PHE:HE2	1.75	0.51
1:A:1086:GLN:NE2	1:A:1118:SER:HB3	2.26	0.51
1:A:1097:ASP:O	1:A:1099:ILE:HG23	2.10	0.51
1:A:189:ASN:OD1	1:A:192:ARG:NH2	2.44	0.51
1:A:942:VAL:HG12	1:A:943:SER:N	2.26	0.51
1:A:1142:GLY:O	1:A:1152:VAL:HA	2.10	0.50
1:A:942:VAL:HG12	1:A:943:SER:H	1.76	0.50
1:A:1027:LEU:HD13	1:A:1083:PHE:HB3	1.94	0.50
1:A:220:LEU:HD22	1:A:956:MET:HG3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1066:ILE:HD12	1:A:1067:GLU:H	1.76	0.50
1:A:54:SER:HB3	1:A:57:ALA:HB3	1.93	0.50
1:A:54:SER:CB	1:A:57:ALA:HB3	2.42	0.49
1:A:1066:ILE:O	1:A:1069:MET:HG2	2.12	0.49
1:A:236:ILE:O	1:A:240:GLU:N	2.43	0.49
1:A:1146:ARG:HH12	1:A:1149:VAL:HB	1.77	0.49
1:A:928:LEU:H	1:A:928:LEU:HD12	1.77	0.49
1:A:8:GLU:HA	1:A:20:VAL:O	2.13	0.49
1:A:208:GLU:HG2	1:A:974:PHE:HE2	1.77	0.49
1:A:1146:ARG:CZ	1:A:1151:ARG:HE	2.25	0.49
1:A:1146:ARG:HH11	1:A:1151:ARG:NH2	2.08	0.48
1:A:909:THR:O	1:A:913:LYS:NZ	2.40	0.48
1:A:1017:LEU:O	1:A:1017:LEU:HD22	2.14	0.48
1:A:224:LEU:HA	1:A:956:MET:CE	2.43	0.48
1:A:1158:GLU:OE1	1:A:1162:LYS:NZ	2.46	0.47
1:A:43:GLY:HA2	1:A:1095:LEU:HD21	1.96	0.47
1:A:986:SER:O	1:A:990:LYS:HG2	2.13	0.47
1:A:201:LYS:HB2	1:A:981:TYR:HE1	1.80	0.47
1:A:245:GLU:O	1:A:249:LYS:HG3	2.15	0.47
1:A:152:ILE:HG13	1:A:1074:LYS:CG	2.41	0.47
1:A:1161:MET:O	1:A:1165:GLU:HG2	2.15	0.46
1:A:247:GLU:HG2	1:A:914:SER:CB	2.44	0.46
1:A:243:ILE:HD11	1:A:917:LEU:C	2.35	0.46
1:A:152:ILE:HG13	1:A:1074:LYS:HD2	1.97	0.46
1:A:1145:MET:C	1:A:1151:ARG:HH12	2.16	0.46
1:A:1003:ASN:O	1:A:1007:LYS:HG2	2.16	0.46
1:A:236:ILE:HG23	1:A:921:LEU:CD2	2.40	0.45
1:A:152:ILE:HG23	1:A:1066:ILE:O	2.16	0.45
1:A:223:ARG:HE	1:A:226:ARG:HH11	1.64	0.45
1:A:935:LEU:O	1:A:939:VAL:HG12	2.16	0.45
1:A:947:GLU:HA	1:A:947:GLU:OE1	2.16	0.45
1:A:1150:SER:C	1:A:1151:ARG:HD2	2.33	0.45
1:A:160:ARG:HG2	1:A:160:ARG:NH1	2.31	0.45
1:A:218:LEU:CD1	1:A:968:MET:HE1	2.46	0.45
1:A:1021:ALA:HA	1:A:1039:LEU:HD12	1.97	0.45
1:A:1146:ARG:N	1:A:1151:ARG:NH2	2.32	0.45
1:A:1104:ASP:O	1:A:1108:VAL:HG13	2.17	0.45
1:A:157:LEU:O	1:A:161:GLN:NE2	2.50	0.45
1:A:193:VAL:O	1:A:197:ILE:HG13	2.17	0.45
1:A:109:ASP:OD1	1:A:111:ARG:HG2	2.17	0.44
1:A:223:ARG:HH21	1:A:226:ARG:CB	2.29	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1057:LYS:HB2	1:A:1063:VAL:HG22	2.00	0.44
1:A:224:LEU:HA	1:A:956:MET:HE3	1.98	0.44
1:A:88:ASN:ND2	1:A:93:PHE:HB2	2.33	0.44
1:A:1146:ARG:HD3	1:A:1151:ARG:NH1	2.13	0.44
1:A:233:LEU:HD13	1:A:940:LYS:HD2	1.99	0.44
1:A:1111:VAL:O	1:A:1115:ILE:HG13	2.18	0.43
1:A:918:ARG:O	1:A:922:GLU:HG2	2.18	0.43
1:A:939:VAL:C	1:A:941:GLU:H	2.20	0.43
1:A:135:ILE:HD13	1:A:1092:PRO:HG2	1.99	0.43
1:A:921:LEU:HD22	1:A:921:LEU:HA	1.64	0.43
1:A:1101:ALA:C	1:A:1102:HIS:HD1	2.22	0.43
1:A:945:ASP:OD1	1:A:948:VAL:HG13	2.19	0.43
1:A:152:ILE:HG13	1:A:1074:LYS:CD	2.49	0.43
1:A:1066:ILE:HA	1:A:1069:MET:CG	2.48	0.43
1:A:1066:ILE:HA	1:A:1069:MET:HG2	2.01	0.43
1:A:205:ASP:HA	1:A:208:GLU:OE1	2.20	0.42
1:A:246:GLY:HA2	1:A:249:LYS:HE3	2.01	0.42
1:A:998:ILE:O	1:A:1002:ILE:HG13	2.19	0.42
1:A:1143:VAL:HA	1:A:1151:ARG:O	2.19	0.42
1:A:68:PHE:CE2	1:A:70:GLY:HA2	2.54	0.42
1:A:945:ASP:O	1:A:948:VAL:HG22	2.19	0.42
1:A:187:GLU:HG2	1:A:187:GLU:H	1.70	0.42
1:A:933:ALA:O	1:A:936:ILE:HG22	2.20	0.42
1:A:14:SER:O	1:A:1150:SER:N	2.49	0.42
1:A:1098:GLU:O	1:A:1099:ILE:HG12	2.19	0.42
1:A:249:LYS:O	1:A:253:ILE:HD12	2.19	0.42
1:A:5:GLU:OE2	1:A:99:GLU:HG3	2.20	0.42
1:A:1146:ARG:O	1:A:1146:ARG:CG	2.67	0.41
1:A:62:ARG:HB3	1:A:62:ARG:HE	1.48	0.41
1:A:926:ARG:HA	1:A:929:LYS:HG3	2.01	0.41
1:A:81:GLU:OE2	1:A:105:ARG:NH1	2.43	0.41
1:A:120:ARG:HA	1:A:120:ARG:HD2	1.78	0.41
1:A:208:GLU:HG2	1:A:974:PHE:CE2	2.56	0.41
1:A:923:GLU:O	1:A:927:GLU:HB2	2.21	0.41
1:A:909:THR:HB	1:A:913:LYS:NZ	2.36	0.41
1:A:218:LEU:HD11	1:A:968:MET:HE1	2.03	0.40
1:A:1066:ILE:CD1	1:A:1067:GLU:OE1	2.70	0.40
1:A:241:SER:HA	1:A:244:ASP:OD2	2.21	0.40
1:A:101:VAL:O	1:A:116:LEU:HA	2.21	0.40
1:A:206:LYS:HA	1:A:206:LYS:HD3	1.91	0.40
1:A:955:ASP:O	1:A:958:ALA:N	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1146:ARG:CZ	1:A:1151:ARG:NE	2.85	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:65:ASP:OD2	1:A:1018:GLU:OE2[1_545]	2.04	0.16

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	513/540 (95%)	474 (92%)	35 (7%)	4 (1%)	22	44

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1099	ILE
1	A	55	ALA
1	A	942	VAL
1	A	962	ALA

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	431/455 (95%)	413 (96%)	18 (4%)	34 62

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	PHE
1	A	64	SER
1	A	131	SER
1	A	223	ARG
1	A	914	SER
1	A	927	GLU
1	A	937	ARG
1	A	956	MET
1	A	997	SER
1	A	1017	LEU
1	A	1046	ASP
1	A	1062	ASP
1	A	1086	GLN
1	A	1087	ARG
1	A	1138	ASP
1	A	1146	ARG
1	A	1147	ASP
1	A	1151	ARG

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

Mol	Chain	Res	Type
1	A	1003	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 carbohydrates 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	517/540 (95%)	0.73	50 (9%) 8 5	35, 59, 113, 143	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	253	ILE	10.2
1	A	254	GLU	8.8
1	A	1169	LYS	8.1
1	A	912	VAL	7.4
1	A	908	ASN	7.1
1	A	910	ILE	7.0
1	A	939	VAL	6.8
1	A	250	ARG	6.7
1	A	907	GLU	6.4
1	A	1164	LEU	6.4
1	A	244	ASP	6.1
1	A	249	LYS	6.1
1	A	940	LYS	5.8
1	A	1168	LYS	5.6
1	A	909	THR	5.5
1	A	1	MET	5.2
1	A	911	LYS	5.0
1	A	922	GLU	4.7
1	A	936	ILE	4.5
1	A	1165	GLU	4.3
1	A	251	ALA	4.0
1	A	937	ARG	3.9
1	A	941	GLU	3.6
1	A	1158	GLU	3.6
1	A	54	SER	3.5
1	A	1166	GLU	3.5
1	A	246	GLY	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	1161	MET	3.1
1	A	1108	VAL	3.1
1	A	1163	ILE	3.0
1	A	252	GLU	2.8
1	A	935	LEU	2.8
1	A	245	GLU	2.8
1	A	913	LYS	2.8
1	A	942	VAL	2.7
1	A	240	GLU	2.5
1	A	944	LEU	2.5
1	A	248	ARG	2.4
1	A	232	LEU	2.4
1	A	950	ARG	2.4
1	A	238	ARG	2.3
1	A	236	ILE	2.3
1	A	929	LYS	2.3
1	A	1160	ALA	2.3
1	A	946	LEU	2.3
1	A	161	GLN	2.2
1	A	239	LEU	2.2
1	A	1162	LYS	2.1
1	A	1087	ARG	2.1
1	A	932	ASP	2.1

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 carbohydrates 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.