Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 09:05 PM GMT

PDB ID	:	4UVK
Title	:	Cohesin subunit Scc3 from Z. rouxii, 88-1035
Authors	:	Roig, M.B.; Nasmyth, K.; Lowe, J.
Deposited on	:	2014-08-06
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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

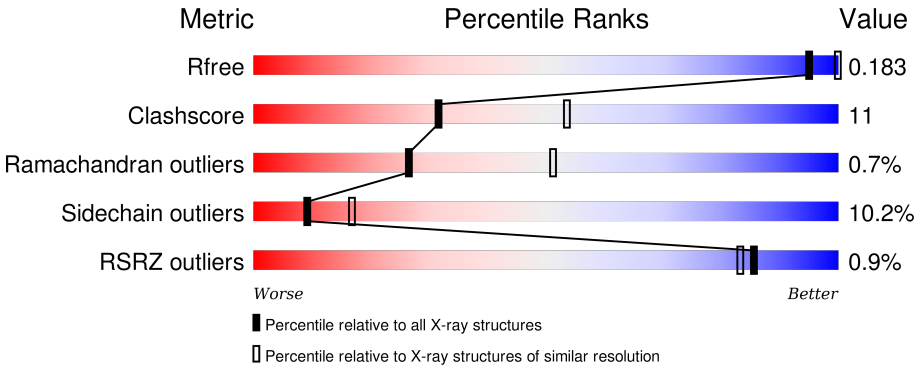
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

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}	91344	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (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 <=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	954	% 67% 22% • 8%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7310 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 ZYRO0D15994P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	876	Total	C	N	O	S	0	0	0
			7101	4603	1160	1320	18			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1036	HIS	-	EXPRESSION TAG	UNP C5DWM3
A	1037	HIS	-	EXPRESSION TAG	UNP C5DWM3
A	1038	HIS	-	EXPRESSION TAG	UNP C5DWM3
A	1039	HIS	-	EXPRESSION TAG	UNP C5DWM3
A	1040	HIS	-	EXPRESSION TAG	UNP C5DWM3
A	1041	HIS	-	EXPRESSION TAG	UNP C5DWM3

- Molecule 2 is water.

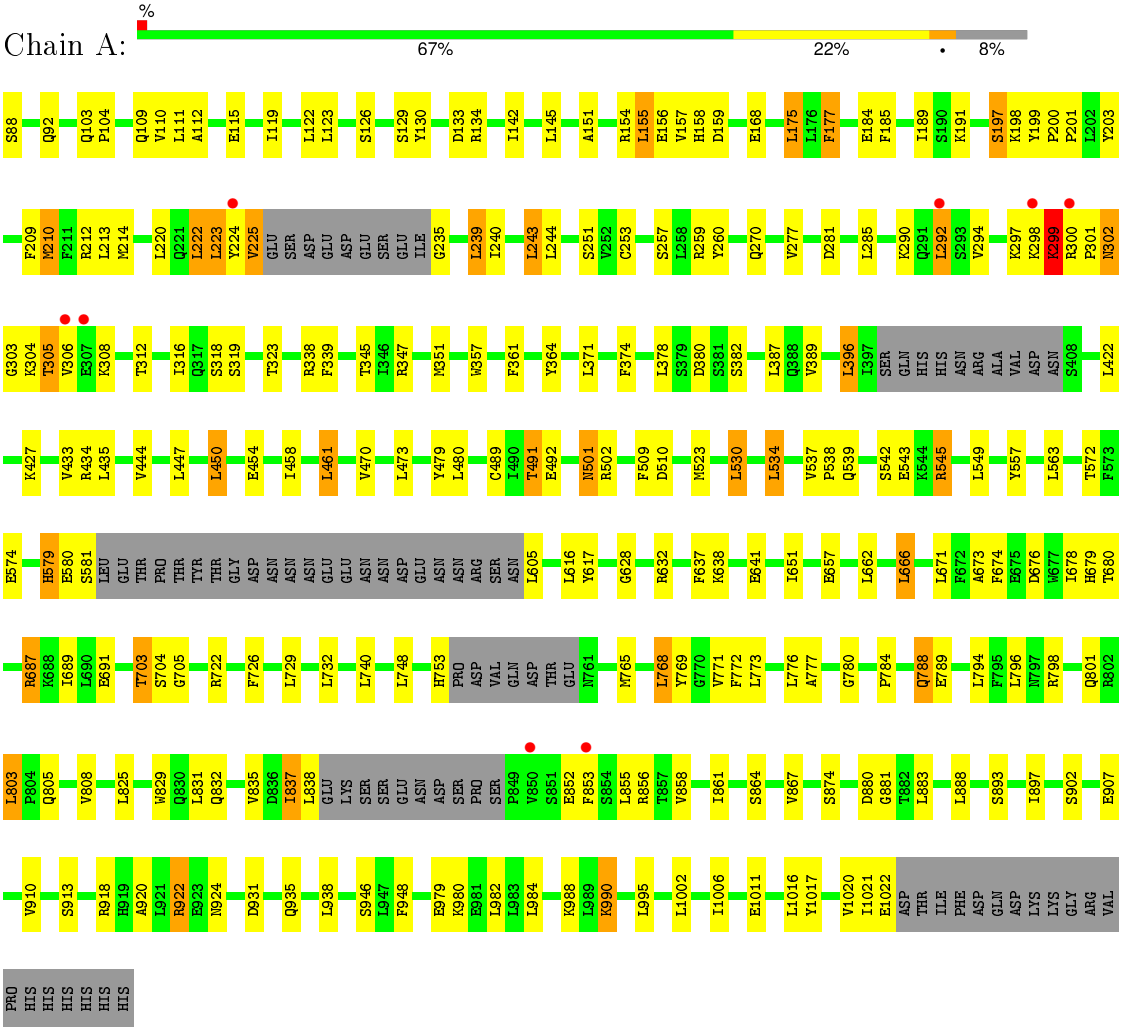
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	209	Total	O	0	0
			209	209		

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 errors 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: ZYRO0D15994P

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	73.37Å 109.18Å 159.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.35 – 2.60 48.35 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.8 (48.35-2.60) 99.6 (48.35-2.60)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.95 (at 2.61Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.187 , 0.245 0.189 , 0.183	Depositor DCC
R_{free} test set	2001 reflections (5.28%)	DCC
Wilson B-factor (Å^2)	39.4	Xtriage
Anisotropy	0.118	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 45.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Outliers	0 of 39910 reflections	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	7310	wwPDB-VP
Average B, all atoms (Å^2)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.73% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	0.74	0/7235	0.77	4/9769 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	435	LEU	CB-CG-CD1	-7.08	98.96	111.00
1	A	396	LEU	CA-CB-CG	6.50	130.24	115.30
1	A	435	LEU	CA-CB-CG	6.09	129.32	115.30
1	A	338	ARG	NE-CZ-NH2	-5.22	117.69	120.30

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	7101	0	7229	153	0
2	A	209	0	0	20	1
All	All	7310	0	7229	153	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 11.

All (153) 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:88:SER:N	2:A:2001:HOH:O	1.88	1.05
1:A:103:GLN:O	1:A:154:ARG:NH1	2.09	0.85
1:A:434:ARG:NH1	2:A:2083:HOH:O	2.09	0.84
1:A:473:LEU:HD21	1:A:632:ARG:CZ	2.09	0.83
1:A:351:MET:HE1	1:A:374:PHE:HA	1.62	0.82
1:A:351:MET:CE	1:A:374:PHE:HA	2.11	0.80
1:A:154:ARG:NE	2:A:2019:HOH:O	1.98	0.79
1:A:235:GLY:N	2:A:2042:HOH:O	2.16	0.78
1:A:530:LEU:HD22	1:A:534:LEU:HD22	1.67	0.76
1:A:220:LEU:HB2	1:A:222:LEU:HD22	1.69	0.74
1:A:298:LYS:O	1:A:300:ARG:N	2.19	0.74
1:A:154:ARG:NH2	2:A:2019:HOH:O	2.20	0.73
1:A:210:MET:CE	1:A:210:MET:HA	2.18	0.73
1:A:225:VAL:O	2:A:2040:HOH:O	2.05	0.72
1:A:302:ASN:O	1:A:302:ASN:ND2	2.22	0.72
1:A:300:ARG:NH1	2:A:2053:HOH:O	2.21	0.71
1:A:154:ARG:CZ	2:A:2019:HOH:O	2.38	0.71
1:A:184:GLU:OE1	1:A:198:LYS:NZ	2.20	0.71
1:A:224:TYR:O	1:A:235:GLY:O	2.10	0.70
1:A:676:ASP:O	1:A:680:THR:HG23	1.92	0.69
1:A:674:PHE:CD1	1:A:726:PHE:CE2	2.82	0.67
1:A:297:LYS:HB3	1:A:298:LYS:HA	1.79	0.65
1:A:210:MET:HE3	1:A:210:MET:HA	1.77	0.63
1:A:765:MET:HE1	1:A:769:TYR:HE2	1.63	0.63
1:A:1006:ILE:O	1:A:1017:TYR:OH	2.13	0.62
1:A:1022:GLU:C	2:A:2209:HOH:O	2.37	0.62
1:A:674:PHE:CZ	1:A:678:ILE:HB	2.35	0.62
1:A:285:LEU:HD21	1:A:316:ILE:HG22	1.82	0.62
1:A:302:ASN:HB3	1:A:305:THR:HB	1.82	0.61
1:A:674:PHE:CD1	1:A:726:PHE:CZ	2.89	0.61
1:A:605:LEU:N	2:A:2117:HOH:O	2.33	0.61
1:A:831:LEU:O	1:A:835:VAL:HG12	2.02	0.59
1:A:579:HIS:O	1:A:581:SER:N	2.31	0.59
1:A:300:ARG:N	1:A:301:PRO:CD	2.65	0.58
1:A:651:ILE:HD12	1:A:689:ILE:HG23	1.85	0.58
1:A:351:MET:HE3	1:A:374:PHE:HD1	1.67	0.58
1:A:509:PHE:O	1:A:510:ASP:HB2	2.03	0.58
1:A:984:LEU:O	1:A:988:LYS:HG2	2.04	0.58
1:A:523:MET:HG2	1:A:617:TYR:CE1	2.39	0.57
1:A:133:ASP:OD1	2:A:2017:HOH:O	2.17	0.57
1:A:557:TYR:CE2	1:A:628:GLY:HA2	2.39	0.57
1:A:239:LEU:HD22	1:A:243:LEU:HD22	1.87	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:224:TYR:CD1	1:A:319:SER:HB2	2.40	0.56
1:A:473:LEU:N	1:A:473:LEU:HD22	2.19	0.56
1:A:210:MET:HE1	1:A:213:LEU:HD12	1.88	0.55
1:A:501:ASN:OD1	1:A:502:ARG:N	2.40	0.55
1:A:557:TYR:CD2	1:A:628:GLY:HA2	2.40	0.55
1:A:351:MET:HE1	1:A:374:PHE:CA	2.35	0.55
1:A:674:PHE:CB	2:A:2136:HOH:O	2.55	0.55
1:A:674:PHE:CE2	1:A:678:ILE:HB	2.41	0.55
1:A:990:LYS:HD2	1:A:1020:VAL:O	2.07	0.55
1:A:220:LEU:CB	1:A:222:LEU:HD22	2.38	0.54
1:A:768:LEU:HD22	1:A:773:LEU:HD22	1.91	0.53
1:A:298:LYS:O	1:A:299:LYS:C	2.46	0.53
1:A:210:MET:HE1	1:A:213:LEU:CD1	2.39	0.52
1:A:768:LEU:HD22	1:A:773:LEU:HD13	1.92	0.52
1:A:224:TYR:O	1:A:225:VAL:C	2.48	0.52
1:A:794:LEU:HD12	1:A:798:ARG:HD2	1.91	0.52
1:A:213:LEU:HD23	1:A:213:LEU:C	2.30	0.52
1:A:458:ILE:HA	1:A:461:LEU:HD22	1.91	0.52
1:A:918:ARG:O	1:A:922:ARG:HG3	2.09	0.52
1:A:458:ILE:O	1:A:461:LEU:HB2	2.10	0.52
1:A:740:LEU:HB3	1:A:776:LEU:HD21	1.92	0.52
1:A:444:VAL:HG12	1:A:450:LEU:HD13	1.91	0.51
1:A:210:MET:HA	1:A:210:MET:HE2	1.92	0.51
1:A:979:GLU:OE2	2:A:2204:HOH:O	2.19	0.51
1:A:537:VAL:HB	1:A:538:PRO:CD	2.40	0.51
1:A:491:THR:HG22	1:A:492:GLU:N	2.25	0.51
1:A:200:PRO:HG2	1:A:201:PRO:HD3	1.94	0.50
1:A:893:SER:O	1:A:897:ILE:HG13	2.12	0.50
1:A:339:PHE:O	1:A:347:ARG:HG2	2.12	0.50
1:A:880:ASP:OD1	1:A:881:GLY:N	2.45	0.49
1:A:371:LEU:HD22	1:A:396:LEU:HD21	1.94	0.49
1:A:637:PHE:CD2	1:A:679:HIS:CE1	3.00	0.49
1:A:461:LEU:HG	1:A:479:TYR:OH	2.13	0.49
1:A:285:LEU:HD21	1:A:316:ILE:CG2	2.43	0.49
1:A:224:TYR:CG	1:A:319:SER:HB2	2.47	0.49
1:A:387:LEU:HD13	1:A:433:VAL:HG22	1.95	0.49
1:A:308:LYS:O	1:A:312:THR:HG23	2.13	0.49
1:A:189:ILE:HG22	1:A:197:SER:HB2	1.94	0.48
1:A:1022:GLU:O	2:A:2209:HOH:O	2.19	0.48
1:A:637:PHE:O	1:A:641:GLU:HG3	2.13	0.48
1:A:780:GLY:HA3	1:A:825:LEU:HD21	1.96	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:361:PHE:CZ	1:A:364:TYR:HB2	2.48	0.48
1:A:537:VAL:HB	1:A:538:PRO:HD2	1.96	0.48
1:A:157:VAL:HG23	1:A:158:HIS:N	2.29	0.48
1:A:224:TYR:CE1	1:A:319:SER:HB2	2.49	0.48
1:A:549:LEU:HD21	1:A:616:LEU:HD12	1.95	0.47
1:A:545:ARG:NH1	2:A:2107:HOH:O	2.43	0.47
1:A:837:ILE:HG22	1:A:838:LEU:HD23	1.97	0.47
1:A:657:GLU:CD	1:A:657:GLU:H	2.17	0.47
1:A:191:LYS:HG3	1:A:203:TYR:CE2	2.50	0.47
1:A:277:VAL:HG23	1:A:323:THR:CG2	2.45	0.46
1:A:253:CYS:O	1:A:259:ARG:HD3	2.15	0.46
1:A:297:LYS:HD2	1:A:299:LYS:H	1.80	0.46
1:A:298:LYS:C	1:A:300:ARG:N	2.66	0.46
1:A:990:LYS:CD	1:A:1020:VAL:O	2.63	0.46
1:A:473:LEU:HD21	1:A:632:ARG:NH1	2.30	0.46
1:A:674:PHE:CD1	1:A:726:PHE:HE2	2.32	0.46
1:A:351:MET:HE3	1:A:389:VAL:HG22	1.97	0.46
1:A:220:LEU:HB2	1:A:222:LEU:CD2	2.42	0.46
1:A:351:MET:HE3	1:A:374:PHE:CD1	2.50	0.46
1:A:214:MET:CE	1:A:240:ILE:HD11	2.45	0.46
1:A:703:THR:HG23	1:A:704:SER:N	2.30	0.46
1:A:662:LEU:HG	1:A:666:LEU:HD22	1.97	0.46
1:A:175:LEU:HA	1:A:175:LEU:HD13	1.77	0.46
1:A:302:ASN:O	1:A:304:LYS:N	2.45	0.45
1:A:572:THR:OG1	2:A:2102:HOH:O	2.10	0.45
1:A:223:LEU:HD23	1:A:240:ILE:HG12	1.98	0.45
1:A:784:PRO:HA	1:A:829:TRP:CE2	2.51	0.45
1:A:549:LEU:CD2	1:A:616:LEU:HD12	2.47	0.45
1:A:938:LEU:HD22	1:A:1002:LEU:HG	1.99	0.45
1:A:674:PHE:N	2:A:2136:HOH:O	2.20	0.45
1:A:920:ALA:O	1:A:924:ASN:ND2	2.45	0.45
1:A:473:LEU:HD21	1:A:632:ARG:NH2	2.31	0.44
1:A:142:ILE:HG21	1:A:155:LEU:HD22	1.99	0.44
1:A:853:PHE:HA	1:A:856:ARG:HB2	1.98	0.44
1:A:260:TYR:HE2	1:A:345:THR:HG22	1.83	0.44
1:A:123:LEU:HD23	1:A:209:PHE:CE1	2.52	0.44
1:A:104:PRO:HB2	1:A:109:GLN:HG2	2.00	0.44
1:A:687:ARG:O	1:A:691:GLU:HG3	2.17	0.43
1:A:480:LEU:HD23	1:A:480:LEU:HA	1.84	0.43
1:A:948:PHE:CZ	1:A:979:GLU:HG3	2.53	0.43
1:A:705:GLY:O	2:A:2138:HOH:O	2.21	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:ALA:HB1	1:A:199:TYR:CD1	2.54	0.43
1:A:673:ALA:O	1:A:676:ASP:HB2	2.19	0.43
1:A:251:SER:O	1:A:259:ARG:CG	2.66	0.43
1:A:92:GLN:NE2	2:A:2002:HOH:O	2.51	0.43
1:A:299:LYS:O	1:A:299:LYS:HG3	2.19	0.43
1:A:130:TYR:CE1	1:A:134:ARG:HG3	2.54	0.43
1:A:351:MET:CE	1:A:374:PHE:CD1	3.02	0.42
1:A:191:LYS:HD2	1:A:203:TYR:CE1	2.54	0.42
1:A:191:LYS:HG3	1:A:203:TYR:CD2	2.55	0.42
1:A:931:ASP:O	1:A:935:GLN:HG2	2.19	0.42
1:A:772:PHE:O	1:A:776:LEU:HG	2.20	0.42
1:A:151:ALA:HA	1:A:177:PHE:CE1	2.55	0.42
1:A:777:ALA:HB2	2:A:2155:HOH:O	2.18	0.42
1:A:726:PHE:HB2	1:A:729:LEU:HD12	2.02	0.42
1:A:156:GLU:O	1:A:159:ASP:HB2	2.20	0.42
1:A:808:VAL:HG12	1:A:808:VAL:O	2.19	0.41
1:A:184:GLU:OE1	1:A:198:LYS:CE	2.68	0.41
1:A:168:GLU:CD	1:A:168:GLU:N	2.74	0.41
1:A:292:LEU:HD22	1:A:292:LEU:O	2.21	0.41
1:A:304:LYS:HG2	1:A:304:LYS:O	2.21	0.41
1:A:347:ARG:NH1	1:A:380:ASP:OD2	2.48	0.41
1:A:422:LEU:HD13	1:A:454:GLU:HG2	2.02	0.41
1:A:297:LYS:CB	1:A:298:LYS:HA	2.49	0.40
1:A:1021:ILE:HA	1:A:1021:ILE:HD13	1.74	0.40
1:A:803:LEU:HG	1:A:888:LEU:HD11	2.03	0.40
1:A:788:GLN:O	1:A:789:GLU:C	2.60	0.40
1:A:796:LEU:O	1:A:801:GLN:HG3	2.21	0.40
1:A:861:ILE:O	1:A:864:SER:HB2	2.22	0.40
1:A:357:TRP:CD1	1:A:357:TRP:N	2.88	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)
2:A:2036:HOH:O	2:A:2052:HOH:O[3_556]	2.19	0.01

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	864/954 (91%)	835 (97%)	23 (3%)	6 (1%)	26 51

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	299	LYS
1	A	580	GLU
1	A	913	SER
1	A	579	HIS
1	A	303	GLY
1	A	837	ILE

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	805/881 (91%)	723 (90%)	82 (10%)	9 17

All (82) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	110	VAL
1	A	111	LEU
1	A	115	GLU
1	A	119	ILE
1	A	122	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	126	SER
1	A	129	SER
1	A	145	LEU
1	A	155	LEU
1	A	175	LEU
1	A	177	PHE
1	A	185	PHE
1	A	197	SER
1	A	210	MET
1	A	212	ARG
1	A	222	LEU
1	A	223	LEU
1	A	225	VAL
1	A	239	LEU
1	A	243	LEU
1	A	244	LEU
1	A	257	SER
1	A	270	GLN
1	A	281	ASP
1	A	290	LYS
1	A	292	LEU
1	A	294	VAL
1	A	299	LYS
1	A	302	ASN
1	A	305	THR
1	A	306	VAL
1	A	318	SER
1	A	378	LEU
1	A	382	SER
1	A	427	LYS
1	A	447	LEU
1	A	450	LEU
1	A	461	LEU
1	A	470	VAL
1	A	489	CYS
1	A	491	THR
1	A	501	ASN
1	A	530	LEU
1	A	534	LEU
1	A	539	GLN
1	A	542	SER
1	A	543	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	545	ARG
1	A	563	LEU
1	A	574	GLU
1	A	638	LYS
1	A	666	LEU
1	A	671	LEU
1	A	687	ARG
1	A	703	THR
1	A	722	ARG
1	A	732	LEU
1	A	748	LEU
1	A	753	HIS
1	A	768	LEU
1	A	771	VAL
1	A	788	GLN
1	A	803	LEU
1	A	805	GLN
1	A	832	GLN
1	A	852	GLU
1	A	855	LEU
1	A	858	VAL
1	A	867	VAL
1	A	874	SER
1	A	883	LEU
1	A	902	SER
1	A	907	GLU
1	A	910	VAL
1	A	922	ARG
1	A	946	SER
1	A	980	LYS
1	A	982	LEU
1	A	990	LYS
1	A	995	LEU
1	A	1011	GLU
1	A	1016	LEU

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	302	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, 95^{th} 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	876/954 (91%)	-0.21	8 (0%)	85	83	21, 38, 73, 116	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	298	LYS	3.9
1	A	292	LEU	2.6
1	A	307	GLU	2.5
1	A	224	TYR	2.5
1	A	306	VAL	2.4
1	A	300	ARG	2.4
1	A	853	PHE	2.1
1	A	850	VAL	2.1

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 carbohydrates 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.