Full wwPDB X-ray Structure Validation Report (i)

Aug 15, 2023 – 10:55 PM EDT

PDB ID	:	1T2P
Title	:	Crystal structure of Sortase A from Staphylococcus aureus
Authors	:	Zong, Y.; Bice, T.W.; Ton-That, H.; Schneewind, O.; Narayana, S.V.
Deposited on	:	2004-04-22
Resolution	:	2.00 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 (1)) were used in the production of this report:

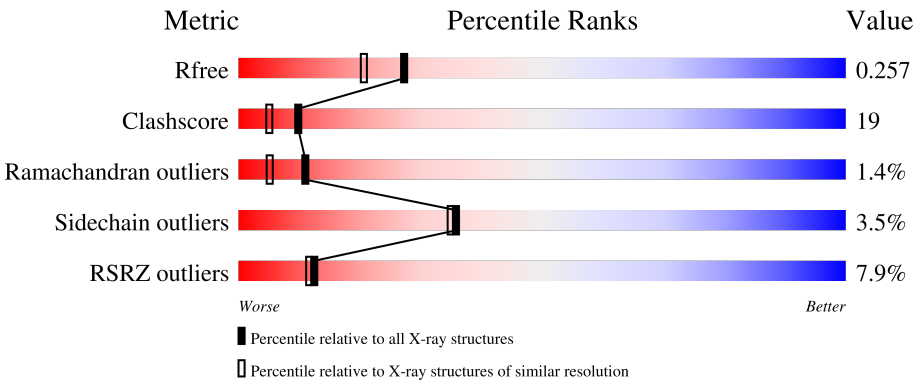
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	146	10% 74% 21% • •
1	B	146	4% 64% 28% • 5%
1	C	146	10% 67% 30% • •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 3759 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 sortase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	145	1155	730	195	227	3	0	0	0
			Total	C	N	O	S			
1	B	138	1091	689	184	215	3	0	0	0
			Total	C	N	O	S			
1	C	145	1151	727	194	227	3	0	0	0

- Molecule 2 is water.

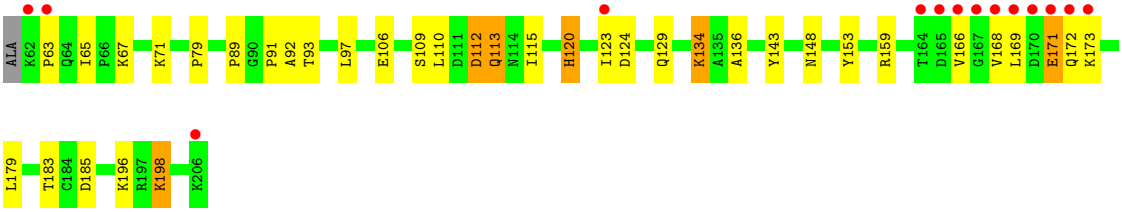
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	107	107	107	0	0
			Total	O		
2	B	146	146	146	0	0
			Total	O		
2	C	109	109	109	0	0

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.

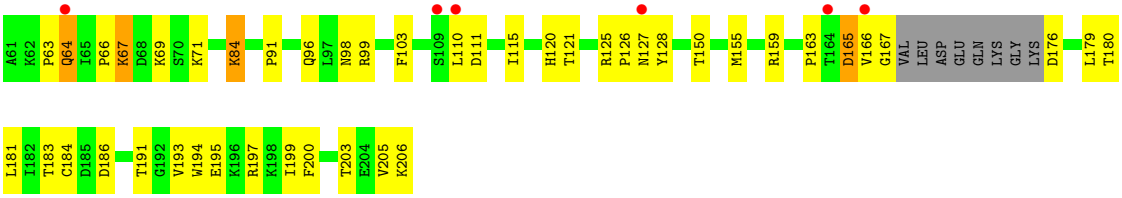
- Molecule 1: sortase

Chain A:



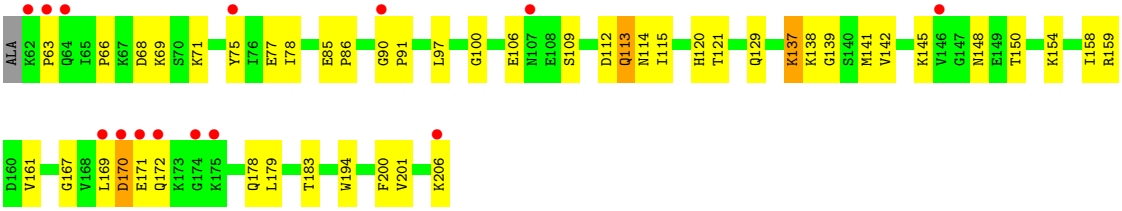
- Molecule 1: sortase

Chain B:



- Molecule 1: sortase

Chain C:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	46.55Å 81.91Å 54.91Å 90.00° 103.75° 90.00°	Depositor
Resolution (Å)	35.52 – 2.00 53.34 – 1.85	Depositor EDS
% Data completeness (in resolution range)	95.2 (35.52-2.00) 97.2 (53.34-1.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.94 (at 1.86Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.209 , 0.257 0.209 , 0.257	Depositor DCC
R_{free} test set	2566 reflections (7.58%)	wwPDB-VP
Wilson B-factor (Å^2)	20.0	Xtriage
Anisotropy	0.041	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 72.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	3759	wwPDB-VP
Average B, all atoms (Å^2)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.18% 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, 0.333 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	0.32	0/1177	0.60	0/1586
1	B	0.31	0/1111	0.59	1/1498 (0.1%)
1	C	0.32	0/1173	0.61	0/1582
All	All	0.32	0/3461	0.60	1/4666 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	63	PRO	N-CA-CB	5.42	109.81	103.30

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	1155	0	1159	42	0
1	B	1091	0	1079	46	0
1	C	1151	0	1148	50	0
2	A	107	0	0	3	0
2	B	146	0	0	9	0
2	C	109	0	0	8	0
All	All	3759	0	3386	130	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 19.

All (130) 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:79:PRO:HG2	1:C:137:LYS:HG2	1.41	1.00
1:C:68:ASP:HB3	1:C:71:LYS:HE2	1.46	0.98
1:B:121:THR:OG1	1:B:183:THR:HB	1.68	0.92
1:C:121:THR:OG1	1:C:183:THR:HB	1.72	0.89
1:C:145:LYS:HE2	1:C:150:THR:OG1	1.76	0.85
1:A:79:PRO:CG	1:C:137:LYS:HG2	2.05	0.85
1:A:91:PRO:HD3	1:A:106:GLU:HB2	1.58	0.84
1:B:96:GLN:HE22	1:B:99:ARG:HH11	1.22	0.84
1:C:172:GLN:H	1:C:178:GLN:HE22	1.27	0.79
1:B:96:GLN:HE22	1:B:99:ARG:NH1	1.83	0.77
1:B:67:LYS:NZ	1:B:67:LYS:HA	1.99	0.77
1:B:64:GLN:NE2	1:B:64:GLN:H	1.82	0.76
1:C:113:GLN:HE22	1:C:172:GLN:HG2	1.51	0.75
1:A:92:ALA:HB1	1:A:97:LEU:HD21	1.70	0.73
1:A:65:ILE:HG13	1:A:110:LEU:HD21	1.71	0.71
1:A:120:HIS:H	1:A:129:GLN:HE21	1.37	0.70
1:C:171:GLU:HB3	1:C:201:VAL:HG21	1.76	0.68
1:B:194:TRP:HB2	1:B:197:ARG:HE	1.59	0.67
1:C:66:PRO:HB3	1:C:71:LYS:HE3	1.76	0.67
1:A:143:TYR:HE1	1:C:137:LYS:HE3	1.59	0.67
1:A:120:HIS:H	1:A:129:GLN:NE2	1.91	0.67
1:B:125:ARG:HD3	2:B:219:HOH:O	1.96	0.66
1:B:71:LYS:HB2	1:B:71:LYS:NZ	2.11	0.66
1:C:172:GLN:H	1:C:178:GLN:NE2	1.94	0.66
1:C:91:PRO:HD3	1:C:106:GLU:HB2	1.78	0.65
1:A:67:LYS:HD2	1:A:67:LYS:H	1.62	0.65
2:A:245:HOH:O	1:B:127:ASN:HB2	1.94	0.65
1:B:121:THR:HG1	1:B:183:THR:HB	1.63	0.64
1:A:79:PRO:HG3	1:C:137:LYS:HE2	1.80	0.63
1:C:137:LYS:HD2	1:C:137:LYS:C	2.19	0.63
1:B:67:LYS:HA	1:B:67:LYS:HZ1	1.63	0.62
1:B:155:MET:HE3	1:B:159:ARG:HD2	1.82	0.62
1:B:180:THR:HG23	1:B:199:ILE:HG23	1.81	0.61
1:C:139:GLY:HA2	1:C:154:LYS:HE2	1.85	0.59
1:C:121:THR:HG1	1:C:183:THR:HB	1.68	0.58
1:B:98:ASN:HB3	2:C:313:HOH:O	2.03	0.58
1:C:114:ASN:HD22	1:C:178:GLN:HG3	1.69	0.58
1:B:71:LYS:HG3	2:B:281:HOH:O	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:VAL:HG12	1:A:168:VAL:HG23	1.86	0.57
1:B:69:LYS:HG3	2:B:315:HOH:O	2.02	0.57
1:A:93:THR:O	1:A:97:LEU:HD23	2.04	0.57
1:A:123:ILE:HG12	2:A:265:HOH:O	2.05	0.57
1:C:121:THR:HG23	2:C:240:HOH:O	2.06	0.56
1:B:64:GLN:NE2	1:B:64:GLN:N	2.54	0.55
1:C:120:HIS:HE1	2:C:208:HOH:O	1.90	0.54
1:B:191:THR:OG1	1:B:193:VAL:HG12	2.08	0.54
1:A:169:LEU:HD23	1:A:169:LEU:H	1.72	0.54
1:B:125:ARG:HD2	1:C:141:MET:HG3	1.90	0.54
1:C:159:ARG:O	1:C:161:VAL:HG13	2.07	0.54
1:B:115:ILE:HG23	1:B:179:LEU:CD1	2.37	0.54
1:B:115:ILE:HG23	1:B:179:LEU:HD13	1.89	0.54
1:A:109:SER:HB2	1:A:112:ASP:OD1	2.08	0.53
1:A:172:GLN:HE21	1:A:172:GLN:HA	1.72	0.53
1:B:110:LEU:HD12	1:B:110:LEU:N	2.24	0.53
1:C:145:LYS:HD3	2:C:243:HOH:O	2.08	0.52
1:A:183:THR:OG1	1:A:198:LYS:HB2	2.09	0.52
1:C:145:LYS:HG2	2:C:266:HOH:O	2.10	0.52
1:B:110:LEU:HD12	1:B:110:LEU:H	1.75	0.52
1:A:65:ILE:HG13	1:A:110:LEU:CD2	2.37	0.52
1:C:115:ILE:HB	1:C:179:LEU:HD12	1.92	0.52
1:B:186:ASP:O	1:B:195:GLU:HG2	2.10	0.51
1:A:143:TYR:HE1	1:C:137:LYS:CE	2.23	0.51
1:C:109:SER:HB3	1:C:112:ASP:OD1	2.11	0.51
1:B:176:ASP:HB3	2:B:298:HOH:O	2.12	0.50
1:B:203:THR:HG23	2:B:234:HOH:O	2.10	0.50
1:B:150:THR:HG22	2:B:329:HOH:O	2.10	0.50
1:B:91:PRO:HD2	2:B:289:HOH:O	2.10	0.49
1:A:171:GLU:O	1:A:171:GLU:HG3	2.11	0.49
1:B:125:ARG:O	1:B:128:TYR:HB3	2.13	0.49
1:C:63:PRO:HD3	1:C:148:ASN:ND2	2.28	0.48
1:B:120:HIS:HD2	1:B:128:TYR:OH	1.96	0.48
1:A:153:TYR:CE2	1:A:179:LEU:HG	2.49	0.48
1:C:63:PRO:HG2	2:C:300:HOH:O	2.14	0.48
1:A:113:GLN:NE2	1:A:172:GLN:HG3	2.29	0.48
1:A:143:TYR:OH	1:C:137:LYS:HG3	2.13	0.48
1:C:69:LYS:HE3	1:C:91:PRO:HD2	1.96	0.48
1:A:97:LEU:N	1:A:97:LEU:HD22	2.29	0.47
1:A:120:HIS:HE1	2:A:212:HOH:O	1.97	0.47
1:A:169:LEU:HD23	1:A:169:LEU:N	2.28	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:134:LYS:H	1:A:134:LYS:HD3	1.79	0.47
1:C:137:LYS:HD3	1:C:138:LYS:O	2.13	0.47
1:B:197:ARG:HD2	2:B:280:HOH:O	2.15	0.47
1:C:171:GLU:HB2	1:C:178:GLN:OE1	2.15	0.47
1:B:84:LYS:O	1:B:84:LYS:HG2	2.16	0.46
1:A:172:GLN:HA	1:A:172:GLN:NE2	2.31	0.45
1:B:103:PHE:HE2	1:B:115:ILE:HD11	1.81	0.45
1:C:121:THR:HG22	2:C:250:HOH:O	2.15	0.45
1:C:137:LYS:HD2	1:C:137:LYS:O	2.17	0.45
1:B:64:GLN:N	1:B:64:GLN:HE21	2.15	0.45
1:B:206:LYS:HD3	2:B:251:HOH:O	2.17	0.45
1:A:110:LEU:HD22	1:A:110:LEU:N	2.32	0.45
1:C:167:GLY:N	1:C:170:ASP:OD1	2.43	0.44
1:C:78:ILE:HG12	1:C:142:VAL:HG22	1.99	0.44
1:A:71:LYS:O	1:A:89:PRO:HD3	2.17	0.44
1:A:168:VAL:HG12	1:A:169:LEU:N	2.32	0.44
1:B:184:CYS:HB3	1:B:194:TRP:CD2	2.53	0.44
1:A:115:ILE:HB	1:A:179:LEU:HD23	2.00	0.44
1:C:139:GLY:HA2	1:C:154:LYS:HG3	1.99	0.44
1:A:67:LYS:HD2	1:A:67:LYS:N	2.31	0.44
1:A:65:ILE:CG1	1:A:110:LEU:HD21	2.45	0.43
1:B:126:PRO:O	1:B:127:ASN:HB3	2.16	0.43
1:C:206:LYS:HA	2:C:313:HOH:O	2.17	0.43
1:C:75:TYR:HE2	1:C:145:LYS:HD2	1.83	0.43
1:A:110:LEU:HD22	1:A:110:LEU:H	1.84	0.43
1:B:71:LYS:HB2	1:B:71:LYS:HZ3	1.80	0.43
1:C:114:ASN:HD22	1:C:178:GLN:CG	2.30	0.43
1:A:134:LYS:HD3	1:A:134:LYS:N	2.34	0.42
1:A:67:LYS:H	1:A:67:LYS:CD	2.31	0.42
1:A:185:ASP:HB3	1:A:196:LYS:HB2	2.00	0.42
1:C:159:ARG:HA	1:C:200:PHE:HA	2.01	0.42
1:C:137:LYS:C	1:C:137:LYS:CD	2.87	0.42
1:B:194:TRP:H	1:B:197:ARG:NH2	2.18	0.42
1:B:165:ASP:O	1:B:167:GLY:N	2.53	0.42
1:C:75:TYR:CD1	1:C:75:TYR:C	2.93	0.42
1:B:66:PRO:HB3	1:B:71:LYS:HD2	2.01	0.41
1:B:155:MET:HE2	1:B:200:PHE:HD2	1.84	0.41
1:B:205:VAL:O	1:B:206:LYS:C	2.59	0.41
1:B:103:PHE:CE2	1:B:115:ILE:HD11	2.54	0.41
1:C:97:LEU:HD12	1:C:194:TRP:CZ2	2.55	0.41
1:C:158:ILE:HG13	1:C:170:ASP:HB3	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:75:TYR:HE1	1:C:77:GLU:HB2	1.85	0.41
1:A:79:PRO:CB	1:C:137:LYS:HG2	2.50	0.41
1:B:193:VAL:HG23	1:B:197:ARG:NH2	2.36	0.41
1:B:163:PRO:HG3	1:B:199:ILE:HD11	2.02	0.41
1:C:85:GLU:HG3	1:C:86:PRO:HD2	2.03	0.41
1:A:113:GLN:CD	1:A:172:GLN:HG3	2.41	0.41
1:C:169:LEU:O	1:C:169:LEU:HG	2.20	0.41
1:A:136:ALA:O	1:A:159:ARG:NE	2.50	0.40
1:C:100:GLY:HA2	1:C:129:GLN:NE2	2.36	0.40
1:A:63:PRO:HG3	1:A:148:ASN:OD1	2.22	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	143/146 (98%)	131 (92%)	10 (7%)	2 (1%)	11	5
1	B	134/146 (92%)	126 (94%)	6 (4%)	2 (2%)	10	4
1	C	143/146 (98%)	134 (94%)	7 (5%)	2 (1%)	11	5
All	All	420/438 (96%)	391 (93%)	23 (6%)	6 (1%)	11	5

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	173	LYS
1	B	166	VAL
1	B	165	ASP
1	C	90	GLY
1	A	171	GLU
1	C	170	ASP

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	128/128 (100%)	122 (95%)	6 (5%)	26	22
1	B	119/128 (93%)	114 (96%)	5 (4%)	30	27
1	C	127/128 (99%)	125 (98%)	2 (2%)	62	67
All	All	374/384 (97%)	361 (96%)	13 (4%)	36	35

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	112	ASP
1	A	113	GLN
1	A	120	HIS
1	A	124	ASP
1	A	134	LYS
1	A	198	LYS
1	B	64	GLN
1	B	67	LYS
1	B	84	LYS
1	B	111	ASP
1	B	181	LEU
1	C	113	GLN
1	C	137	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	113	GLN
1	A	120	HIS
1	A	129	GLN
1	A	172	GLN
1	A	178	GLN
1	B	64	GLN
1	B	96	GLN
1	B	98	ASN

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Mol	Chain	Res	Type
1	B	107	ASN
1	B	113	GLN
1	B	120	HIS
1	C	113	GLN
1	C	114	ASN
1	C	120	HIS
1	C	148	ASN
1	C	178	GLN

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 (i)

6.1 Protein, DNA and RNA chains (i)

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	145/146 (99%)	0.27	14 (9%)	7	7	11, 20, 104, 131	0
1	B	138/146 (94%)	-0.02	6 (4%)	35	34	10, 21, 69, 88	0
1	C	145/146 (99%)	0.18	14 (9%)	7	7	7, 19, 84, 137	0
All	All	428/438 (97%)	0.15	34 (7%)	12	11	7, 20, 84, 137	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	167	GLY	12.3
1	C	172	GLN	7.6
1	C	90	GLY	7.1
1	C	171	GLU	6.8
1	A	168	VAL	6.7
1	A	169	LEU	6.6
1	A	170	ASP	5.9
1	C	107	ASN	5.7
1	A	166	VAL	5.5
1	C	206	LYS	4.8
1	A	206	LYS	4.8
1	C	75	TYR	4.7
1	A	173	LYS	4.6
1	A	172	GLN	4.4
1	B	166	VAL	4.3
1	A	165	ASP	3.9
1	A	171	GLU	3.9
1	A	164	THR	3.8
1	A	62	LYS	3.8
1	C	175	LYS	3.6
1	C	169	LEU	3.5
1	B	127	ASN	3.3
1	B	64	GLN	3.1

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Mol	Chain	Res	Type	RSRZ
1	C	170	ASP	3.1
1	B	110	LEU	2.9
1	B	164	THR	2.7
1	C	63	PRO	2.6
1	A	63	PRO	2.4
1	A	123	ILE	2.4
1	C	146	VAL	2.2
1	C	174	GLY	2.1
1	B	109	SER	2.1
1	C	62	LYS	2.1
1	C	64	GLN	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 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.