Full wwPDB X-ray Structure Validation Report (i)

Oct 24, 2023 – 03:50 PM EDT

PDB ID	:	3AH6
Title	:	Remarkable improvement of the heat stability of CutA1 from E.coli by rational protein designing
Authors	:	Matsuura, Y.; Tanaka, T.; Bagautdinov, B.; Kunishima, N.; Yutani, K.
Deposited on	:	2010-04-15
Resolution	:	2.40 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:

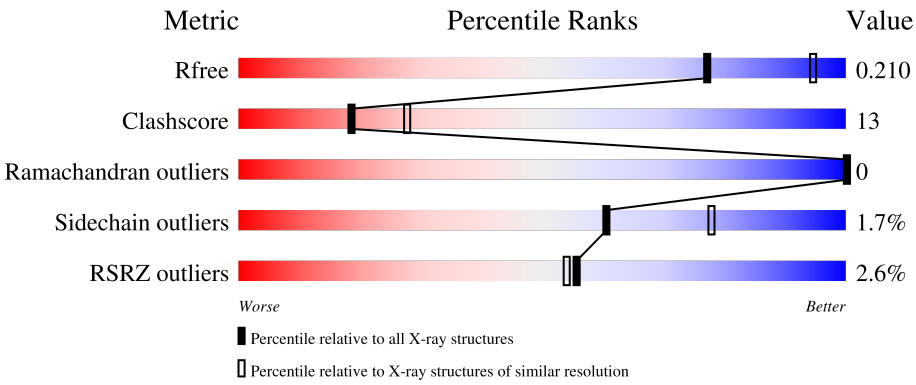
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	112	75% 16% • 8%
1	B	112	4% 73% 20% 7%
1	C	112	73% 21% 6%
1	D	112	4% 75% 17% • 7%
1	E	112	3% 70% 21% • 8%

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Mol	Chain	Length	Quality of chain
1	F	112	4% 58% 33% 8%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 5004 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 Divalent-cation tolerance protein cutA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	103	792	509	125	154	4	0	0	0
1	B	104	800	513	127	156	4	0	0	0
1	C	105	806	516	128	158	4	0	0	0
1	D	104	804	515	129	156	4	0	0	0
1	E	103	792	509	125	154	4	0	0	0
1	F	103	792	509	125	154	4	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	11	VAL	SER	engineered mutation	UNP P69488
B	11	VAL	SER	engineered mutation	UNP P69488
C	11	VAL	SER	engineered mutation	UNP P69488
D	11	VAL	SER	engineered mutation	UNP P69488
E	11	VAL	SER	engineered mutation	UNP P69488
F	11	VAL	SER	engineered mutation	UNP P69488

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	38	38	38	0	0
2	B	44	44	44	0	0
2	C	43	43	43	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	D	40	40	40	0	0
			Total	O		
2	E	32	32	32	0	0
			Total	O		
2	F	21	21	21	0	0

3 Residue-property plots (i)

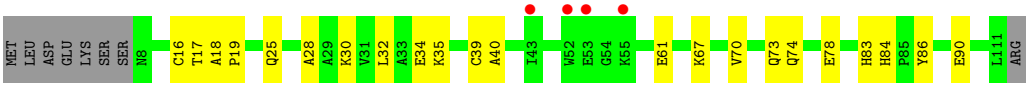
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Divalent-cation tolerance protein cutA

Chain A:

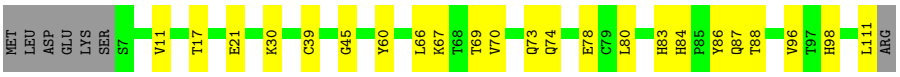
• Molecule 1: Divalent-cation tolerance protein cutA

Chain B:



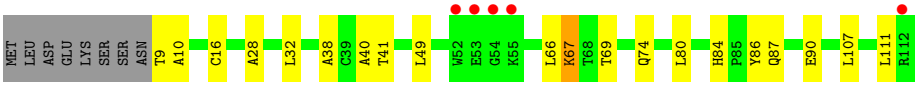
• Molecule 1: Divalent-cation tolerance protein cutA

Chain C:



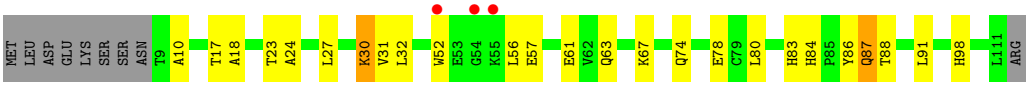
• Molecule 1: Divalent-cation tolerance protein cutA

Chain D:



• Molecule 1: Divalent-cation tolerance protein cutA

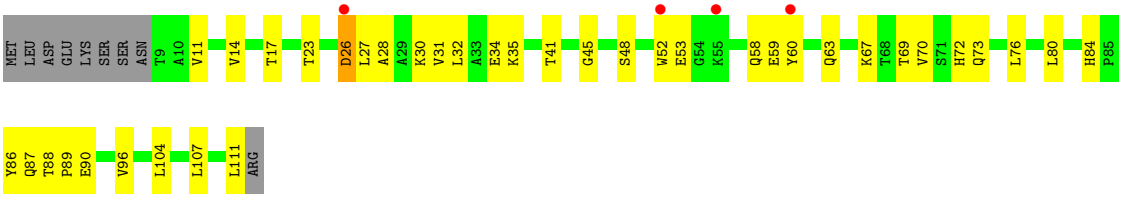
Chain E:



• Molecule 1: Divalent-cation tolerance protein cutA

Chain F:

4%
58%
33%
8%


MET LEU ASP GLU LYS SER SER ASN T9 A10 V11 V14 T17 T23 D26 L27 A28 A29 K30 V31 L32 A33 E34 K35 T41 G45 S48 W52 E53 G54 K55 Q58 E59 Y60 Q63 K67 T68 T69 V70 S71 H72 Q73 L76 L80 H84 P85

Y86 Q87 T88 P89 E90 V96 L104 L107 L111 ARG

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	62.24Å 96.87Å 106.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.23 – 2.40 38.22 – 2.40	Depositor EDS
% Data completeness (in resolution range)	100.0 (38.23-2.40) 99.9 (38.22-2.40)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.98 (at 2.39Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.202 , 0.251 0.203 , 0.210	Depositor DCC
R_{free} test set	1316 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å^2)	35.1	Xtriage
Anisotropy	0.075	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 45.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	5004	wwPDB-VP
Average B, all atoms (Å^2)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.79% 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	0.34	0/809	0.58	0/1110
1	B	0.33	0/817	0.56	0/1121
1	C	0.34	0/823	0.57	0/1129
1	D	0.32	0/821	0.58	0/1124
1	E	0.36	0/809	0.61	0/1110
1	F	0.35	0/809	0.58	0/1110
All	All	0.34	0/4888	0.58	0/6704

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 (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	792	0	793	15	0
1	B	800	0	799	17	0
1	C	806	0	804	19	0
1	D	804	0	806	20	0
1	E	792	0	793	28	0
1	F	792	0	793	36	0
2	A	38	0	0	2	0
2	B	44	0	0	0	0
2	C	43	0	0	2	0
2	D	40	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	32	0	0	1	0
2	F	21	0	0	1	0
All	All	5004	0	4788	126	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 13.

All (126) 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:E:30:LYS:HB2	1:E:30:LYS:NZ	1.83	0.93
1:E:84:HIS:HD2	1:E:86:TYR:H	1.15	0.91
1:B:84:HIS:HD2	1:B:86:TYR:H	1.19	0.88
1:C:84:HIS:HD2	1:C:86:TYR:H	1.22	0.85
1:D:84:HIS:HD2	1:D:86:TYR:H	1.25	0.84
1:F:84:HIS:HD2	1:F:86:TYR:H	1.29	0.78
1:E:30:LYS:HB2	1:E:30:LYS:HZ2	1.50	0.75
1:F:87:GLN:NE2	1:F:88:THR:H	1.86	0.74
1:B:74:GLN:O	1:B:78:GLU:HG3	1.88	0.73
1:E:87:GLN:H	1:E:87:GLN:CD	1.93	0.72
1:E:87:GLN:H	1:E:87:GLN:NE2	1.89	0.71
1:F:23:THR:O	1:F:27:LEU:HB2	1.93	0.69
1:F:72:HIS:HD2	1:F:111:LEU:HD13	1.59	0.68
1:A:21:GLU:O	1:A:25:GLN:HG3	1.94	0.67
1:F:87:GLN:HE21	1:F:88:THR:H	1.40	0.67
1:E:84:HIS:CD2	1:E:86:TYR:H	2.07	0.67
1:C:87:GLN:H	1:C:87:GLN:CD	1.98	0.67
1:D:38:ALA:HB3	1:D:107:LEU:HD23	1.77	0.67
1:F:69:THR:HG23	1:F:111:LEU:HD12	1.77	0.67
1:D:9:THR:HA	2:D:211:HOH:O	1.95	0.66
1:D:87:GLN:H	1:D:87:GLN:CD	2.01	0.64
1:E:87:GLN:HG2	1:E:88:THR:H	1.63	0.62
1:A:23:THR:O	1:A:27:LEU:HB2	1.99	0.62
1:A:9:THR:N	2:A:224:HOH:O	2.33	0.62
1:A:84:HIS:CD2	1:A:89:PRO:HB3	2.35	0.61
1:E:30:LYS:HB2	1:E:30:LYS:HZ3	1.64	0.61
1:F:45:GLY:HA2	1:F:60:TYR:HE1	1.65	0.61
1:F:26:ASP:O	1:F:30:LYS:HG3	2.00	0.60
1:E:30:LYS:HE2	1:E:83:HIS:CE1	2.37	0.60
1:F:28:ALA:O	1:F:32:LEU:HB2	2.03	0.59
1:B:30:LYS:HE2	1:B:83:HIS:CE1	2.38	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:84:HIS:CD2	1:F:86:TYR:H	2.18	0.58
1:A:70:VAL:O	1:A:73:GLN:HG2	2.02	0.58
1:F:72:HIS:CD2	1:F:111:LEU:HD13	2.38	0.58
1:D:67:LYS:N	1:D:67:LYS:HD3	2.18	0.57
1:B:30:LYS:HE2	1:B:83:HIS:NE2	2.19	0.57
1:F:26:ASP:OD1	1:F:30:LYS:HE2	2.04	0.57
1:B:18:ALA:O	1:B:61:GLU:HB3	2.05	0.57
1:E:87:GLN:NE2	1:E:87:GLN:N	2.53	0.57
1:B:70:VAL:O	1:B:73:GLN:HG2	2.04	0.56
1:D:84:HIS:CD2	1:D:86:TYR:H	2.16	0.56
1:A:30:LYS:HE2	1:A:83:HIS:CE1	2.41	0.55
1:B:28:ALA:O	1:B:32:LEU:HD13	2.06	0.55
1:E:17:THR:HG21	1:E:61:GLU:OE1	2.06	0.55
1:A:67:LYS:HD2	2:A:221:HOH:O	2.07	0.55
1:B:17:THR:O	1:B:84:HIS:HE1	1.89	0.55
1:E:56:LEU:CD1	1:F:32:LEU:HD23	2.38	0.54
1:F:70:VAL:O	1:F:73:GLN:HG2	2.08	0.54
1:A:98:HIS:HB2	2:C:202:HOH:O	2.08	0.53
1:A:72:HIS:HD2	1:A:111:LEU:HD13	1.72	0.53
1:C:70:VAL:O	1:C:73:GLN:HG2	2.09	0.53
1:E:17:THR:O	1:E:84:HIS:HE1	1.93	0.52
1:C:30:LYS:HE2	1:C:83:HIS:NE2	2.25	0.52
1:C:69:THR:HG23	1:C:111:LEU:HD12	1.92	0.51
1:D:28:ALA:O	1:D:32:LEU:HG	2.11	0.51
1:D:49:LEU:HD12	1:D:49:LEU:N	2.26	0.50
1:F:11:VAL:HG23	1:F:96:VAL:HB	1.94	0.50
1:D:67:LYS:HE2	1:F:90:GLU:OE2	2.12	0.49
1:C:84:HIS:CD2	1:C:86:TYR:H	2.14	0.49
1:D:84:HIS:CD2	1:D:86:TYR:HB2	2.47	0.49
1:F:17:THR:O	1:F:84:HIS:HE1	1.95	0.49
1:D:84:HIS:HD2	1:D:86:TYR:N	2.03	0.49
1:C:39:CYS:SG	1:C:67:LYS:HE2	2.53	0.49
1:F:84:HIS:HD2	1:F:86:TYR:N	2.07	0.48
1:F:72:HIS:CD2	1:F:111:LEU:HB3	2.50	0.47
1:C:69:THR:HG23	1:C:111:LEU:CD1	2.44	0.47
1:A:72:HIS:CD2	1:A:111:LEU:HD13	2.48	0.47
1:F:84:HIS:CD2	1:F:86:TYR:HB2	2.50	0.47
1:A:17:THR:O	1:A:84:HIS:HE1	1.98	0.47
1:C:11:VAL:HG23	1:C:96:VAL:HG21	1.96	0.47
1:C:17:THR:O	1:C:84:HIS:HE1	1.98	0.47
1:E:10:ALA:HB2	2:E:212:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:87:GLN:HE21	1:F:88:THR:N	2.11	0.46
1:C:66:LEU:HD12	1:C:66:LEU:N	2.30	0.46
1:A:84:HIS:HD2	1:A:86:TYR:H	1.63	0.46
1:F:58:GLN:O	1:F:59:GLU:HG3	2.16	0.46
1:E:84:HIS:CD2	1:E:86:TYR:HB2	2.50	0.46
1:C:80:LEU:O	1:C:84:HIS:HB2	2.16	0.46
1:B:84:HIS:CD2	1:B:86:TYR:H	2.11	0.45
1:B:74:GLN:OE1	1:C:98:HIS:HE1	1.99	0.45
1:D:40:ALA:O	1:F:48:SER:HA	2.17	0.45
1:D:16:CYS:HA	1:D:90:GLU:O	2.17	0.45
1:A:30:LYS:HE2	1:A:83:HIS:NE2	2.31	0.45
1:D:41:THR:HG21	1:F:63:GLN:HE21	1.82	0.45
1:F:45:GLY:HA2	1:F:60:TYR:CE1	2.49	0.45
1:F:31:VAL:HG11	1:F:80:LEU:CD2	2.47	0.45
1:E:31:VAL:HG11	1:E:80:LEU:CD2	2.47	0.45
1:D:74:GLN:HE22	1:E:98:HIS:CE1	2.36	0.44
1:D:80:LEU:O	1:D:84:HIS:HB2	2.18	0.44
1:B:16:CYS:HA	1:B:90:GLU:O	2.18	0.43
1:C:98:HIS:HD2	2:C:178:HOH:O	2.01	0.43
1:D:69:THR:HG23	1:D:111:LEU:HD12	2.00	0.43
1:E:17:THR:CG2	1:E:61:GLU:HB3	2.49	0.43
1:E:87:GLN:HG2	1:E:88:THR:N	2.31	0.43
1:E:84:HIS:HD2	1:E:86:TYR:N	1.99	0.43
1:B:34:GLU:O	1:B:35:LYS:HB2	2.19	0.43
1:A:80:LEU:O	1:A:84:HIS:HB2	2.18	0.43
1:F:84:HIS:CD2	1:F:89:PRO:HB3	2.54	0.43
1:C:45:GLY:HA2	1:C:60:TYR:HE1	1.83	0.42
1:F:31:VAL:HG11	1:F:80:LEU:HG	2.01	0.42
1:C:87:GLN:HG2	1:C:88:THR:H	1.83	0.42
1:F:11:VAL:HG11	1:F:107:LEU:HD23	2.01	0.42
1:F:52:TRP:CD1	1:F:53:GLU:HG2	2.54	0.42
1:C:87:GLN:HG2	1:C:88:THR:N	2.34	0.42
1:F:30:LYS:O	1:F:34:GLU:HG2	2.18	0.42
1:B:19:PRO:HB3	1:B:86:TYR:OH	2.19	0.42
1:B:84:HIS:CD2	1:B:86:TYR:HB2	2.55	0.42
1:E:30:LYS:HG3	1:E:31:VAL:N	2.35	0.42
1:E:18:ALA:HB3	1:E:24:ALA:HB2	2.02	0.42
1:F:34:GLU:O	1:F:35:LYS:HB2	2.20	0.41
1:F:45:GLY:CA	1:F:60:TYR:HE1	2.32	0.41
1:B:25:GLN:HE22	1:D:10:ALA:CB	2.33	0.41
1:E:23:THR:O	1:E:27:LEU:HB2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:11:VAL:HG22	2:F:123:HOH:O	2.19	0.41
1:E:91:LEU:HD23	1:E:91:LEU:HA	1.95	0.41
1:C:21:GLU:OE2	1:C:60:TYR:OH	2.25	0.41
1:C:74:GLN:O	1:C:78:GLU:HG3	2.20	0.41
1:E:52:TRP:HB3	1:E:57:GLU:HG3	2.03	0.41
1:E:74:GLN:O	1:E:78:GLU:HG3	2.21	0.41
1:F:14:VAL:HG21	1:F:76:LEU:HD23	2.02	0.41
1:B:39:CYS:HB3	1:B:67:LYS:HE2	2.03	0.41
1:E:63:GLN:HE21	1:F:41:THR:HG21	1.85	0.41
1:D:67:LYS:HB3	1:D:107:LEU:HD21	2.03	0.41
1:A:48:SER:HA	1:B:40:ALA:O	2.21	0.40
1:E:80:LEU:O	1:E:84:HIS:HB2	2.21	0.40
1:D:40:ALA:HB2	1:D:66:LEU:HD23	2.02	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	101/112 (90%)	100 (99%)	1 (1%)	0	100	100
1	B	102/112 (91%)	102 (100%)	0	0	100	100
1	C	103/112 (92%)	103 (100%)	0	0	100	100
1	D	102/112 (91%)	102 (100%)	0	0	100	100
1	E	101/112 (90%)	98 (97%)	3 (3%)	0	100	100
1	F	101/112 (90%)	99 (98%)	2 (2%)	0	100	100
All	All	610/672 (91%)	604 (99%)	6 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	87/96 (91%)	86 (99%)	1 (1%)	73	87
1	B	88/96 (92%)	88 (100%)	0	100	100
1	C	89/96 (93%)	89 (100%)	0	100	100
1	D	88/96 (92%)	87 (99%)	1 (1%)	73	87
1	E	87/96 (91%)	83 (95%)	4 (5%)	27	43
1	F	87/96 (91%)	84 (97%)	3 (3%)	37	56
All	All	526/576 (91%)	517 (98%)	9 (2%)	60	78

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	LYS
1	D	67	LYS
1	E	30	LYS
1	E	32	LEU
1	E	67	LYS
1	E	87	GLN
1	F	26	ASP
1	F	67	LYS
1	F	104	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	HIS
1	A	73	GLN
1	A	74	GLN
1	A	84	HIS
1	B	25	GLN
1	B	84	HIS
1	B	98	HIS
1	C	84	HIS

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Mol	Chain	Res	Type
1	C	98	HIS
1	D	25	GLN
1	D	74	GLN
1	D	84	HIS
1	D	98	HIS
1	D	108	ASN
1	E	63	GLN
1	E	84	HIS
1	E	87	GLN
1	E	98	HIS
1	F	25	GLN
1	F	63	GLN
1	F	72	HIS
1	F	84	HIS
1	F	87	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	103/112 (91%)	-0.07	0	100	100	20, 32, 42, 55	0
1	B	104/112 (92%)	0.00	4 (3%)	40	39	19, 29, 51, 66	0
1	C	105/112 (93%)	-0.20	0	100	100	21, 31, 41, 48	0
1	D	104/112 (92%)	-0.15	5 (4%)	30	29	22, 32, 49, 65	0
1	E	103/112 (91%)	0.08	3 (2%)	51	50	23, 35, 57, 61	0
1	F	103/112 (91%)	0.25	4 (3%)	39	38	29, 42, 55, 61	0
All	All	622/672 (92%)	-0.02	16 (2%)	56	54	19, 33, 54, 66	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	52	TRP	5.8
1	B	52	TRP	4.4
1	D	112	ARG	3.8
1	E	55	LYS	3.3
1	E	54	GLY	3.3
1	F	52	TRP	3.2
1	F	55	LYS	3.0
1	D	53	GLU	3.0
1	E	52	TRP	2.9
1	D	54	GLY	2.6
1	F	26	ASP	2.6
1	F	60	TYR	2.6
1	B	53	GLU	2.3
1	D	55	LYS	2.2
1	B	55	LYS	2.1
1	B	43	ILE	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.