



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 24, 2023 – 09:43 AM EDT

PDB ID : 2ZFH
Title : Crystal structure of putative CutA1 from Homo sapiens at 2.05Å resolution
Authors : Bagautdinov, B.; Yutani, K.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2008-01-07
Resolution : 2.05 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 ⓘ 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 ⓘ](#)) 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

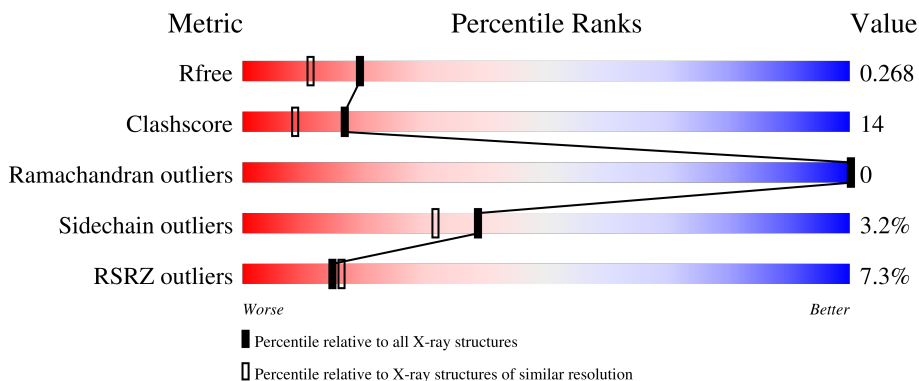
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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.



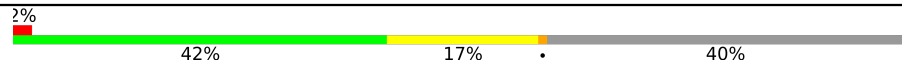
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1692 (2.04-2.04)
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)
RSRZ outliers	127900	1672 (2.04-2.04)

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 $\leq 5\%$. 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	179	 4% 40% 21% 39%
1	B	179	 3% 43% 16% 40%
1	C	179	 3% 42% 16% 41%
1	D	179	 9% 39% 20% 40%
1	E	179	 4% 47% 13% 40%

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Mol	Chain	Length	Quality of chain
1	F	179	 <p>A horizontal bar chart representing the quality of the chain. The bar is divided into four segments: a small red segment at the beginning labeled '2%', followed by a green segment labeled '42%', a yellow segment labeled '17%', and a grey segment at the end labeled '40%'. A small black dot is located on the grey segment.</p>

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 5279 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 CutA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	109	851	548	139	160	4	0	0	0
1	B	107	836	540	137	155	4	0	0	0
1	C	106	832	538	136	154	4	0	0	0
1	D	107	836	540	137	155	4	0	0	0
1	E	107	836	540	137	155	4	0	0	0
1	F	107	836	540	137	155	4	0	0	0

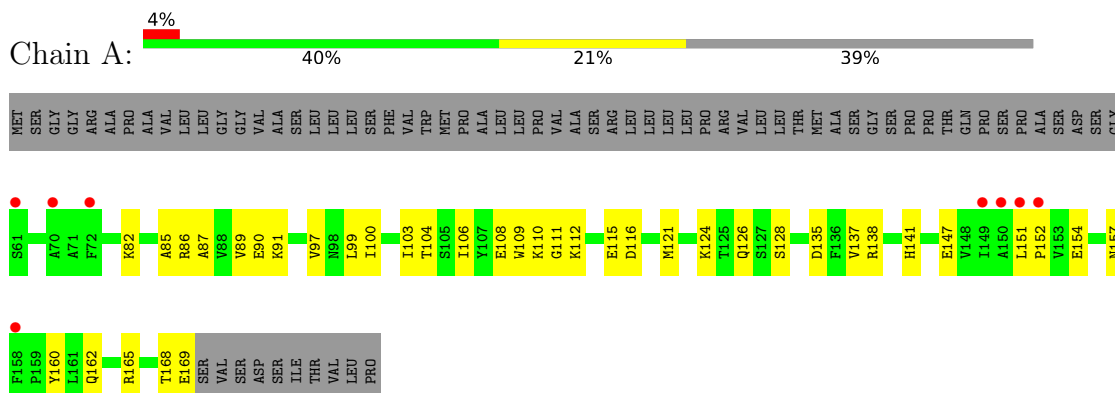
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	35	Total	O	0	0
			35	35		
2	B	37	Total	O	0	0
			37	37		
2	C	50	Total	O	0	0
			50	50		
2	D	48	Total	O	0	0
			48	48		
2	E	44	Total	O	0	0
			44	44		
2	F	38	Total	O	0	0
			38	38		

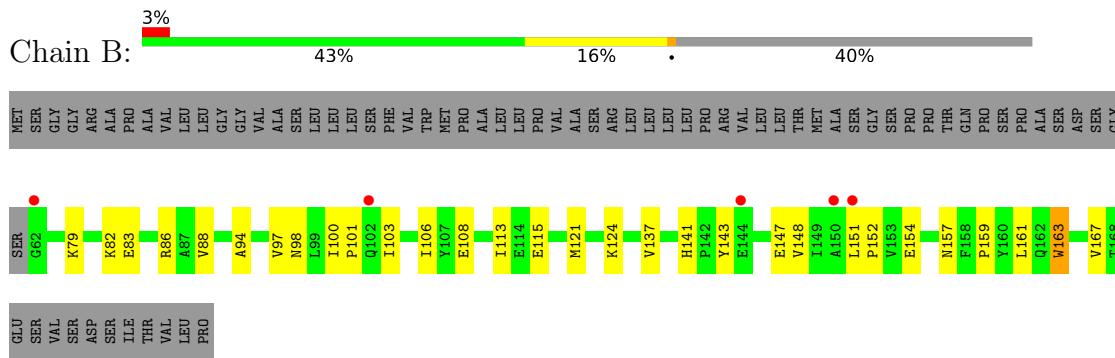
3 Residue-property plots i

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.

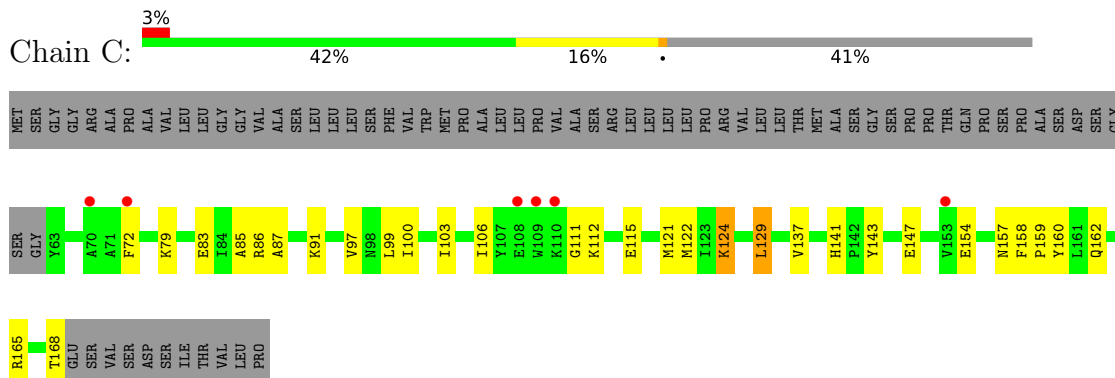
- Molecule 1: CutA



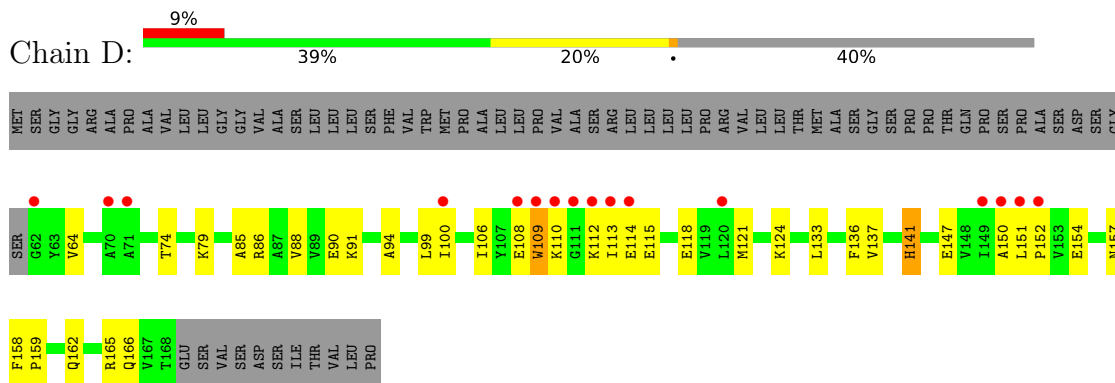
- Molecule 1: CutA



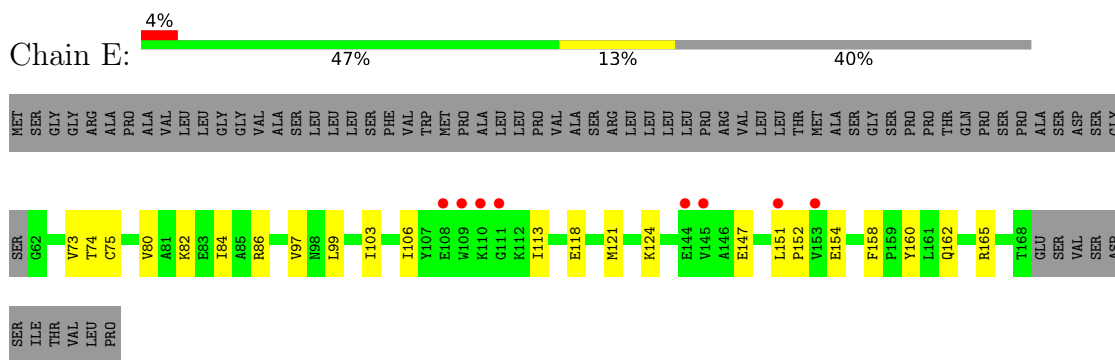
- Molecule 1: CutA



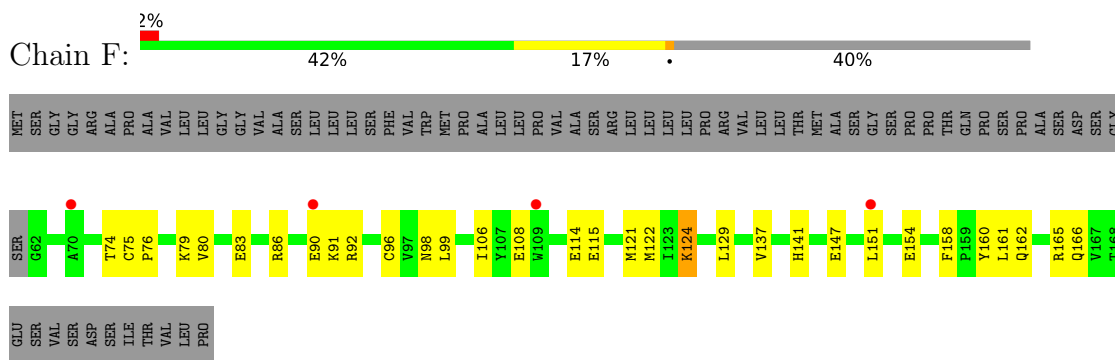
• Molecule 1: CutA



• Molecule 1: CutA



• Molecule 1: CutA



4 Data and refinement statistics i

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	68.69Å 88.84Å 125.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.81 – 2.05 37.81 – 2.06	Depositor EDS
% Data completeness (in resolution range)	94.7 (37.81-2.05) 94.8 (37.81-2.06)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.86 (at 2.06Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.221 , 0.265 0.223 , 0.268	Depositor DCC
R_{free} test set	2291 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	35.9	Xtrriage
Anisotropy	0.906	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 53.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5279	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.80% 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	0.35	0/869	0.57	0/1185
1	B	0.35	0/854	0.60	0/1165
1	C	0.38	0/850	0.59	0/1160
1	D	0.40	0/854	0.63	0/1165
1	E	0.35	0/854	0.57	0/1165
1	F	0.37	0/854	0.58	0/1165
All	All	0.37	0/5135	0.59	0/7005

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	851	0	863	30	0
1	B	836	0	852	24	0
1	C	832	0	849	26	0
1	D	836	0	852	39	0
1	E	836	0	852	24	0
1	F	836	0	852	28	0
2	A	35	0	0	0	0
2	B	37	0	0	0	0
2	C	50	0	0	0	0
2	D	48	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	44	0	0	2	0
2	F	38	0	0	1	0
All	All	5279	0	5120	143	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 14.

All (143) 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:D:108:GLU:HG3	1:D:113:ILE:HD13	1.42	0.97
1:D:162:GLN:HG3	1:D:165:ARG:NH2	1.92	0.84
1:C:122:MET:HG2	1:C:124:LYS:HZ3	1.46	0.80
1:B:100:ILE:HG22	1:B:103:ILE:HD11	1.62	0.80
1:B:106:ILE:HG12	1:B:115:GLU:HG2	1.69	0.73
1:E:160:TYR:CZ	1:F:147:GLU:HG3	2.24	0.72
1:A:111:GLY:O	1:A:112:LYS:HD2	1.89	0.72
1:C:79:LYS:O	1:C:83:GLU:HG3	1.90	0.70
1:E:162:GLN:HG3	1:E:165:ARG:NH2	2.06	0.70
1:B:137:VAL:O	1:B:141:HIS:HB2	1.93	0.69
1:B:79:LYS:O	1:B:83:GLU:HG3	1.94	0.68
1:C:162:GLN:HG3	1:C:165:ARG:HH21	1.57	0.68
1:A:82:LYS:O	1:A:86:ARG:HG3	1.94	0.68
1:A:160:TYR:CZ	1:B:147:GLU:HG3	2.29	0.67
1:A:106:ILE:HG12	1:A:115:GLU:HG2	1.75	0.67
1:C:106:ILE:HG12	1:C:115:GLU:HG2	1.77	0.67
1:A:162:GLN:HG3	1:A:165:ARG:NH2	2.12	0.64
1:D:109:TRP:HD1	1:D:114:GLU:OE2	1.81	0.64
1:F:162:GLN:O	1:F:166:GLN:HG3	1.96	0.63
1:E:74:THR:HG22	1:E:147:GLU:HB3	1.80	0.63
1:F:79:LYS:O	1:F:83:GLU:HG3	1.99	0.63
1:F:76:PRO:HD2	1:F:80:VAL:HG21	1.81	0.63
1:D:113:ILE:HD11	1:F:90:GLU:HG3	1.81	0.62
1:D:158:PHE:HB3	2:D:181:HOH:O	1.99	0.61
1:F:165:ARG:HG3	1:F:165:ARG:HH11	1.65	0.61
1:F:106:ILE:HG12	1:F:115:GLU:HG2	1.83	0.60
1:C:122:MET:HG2	1:C:124:LYS:NZ	2.15	0.60
1:F:108:GLU:HG3	2:F:213:HOH:O	2.00	0.60
1:D:108:GLU:HG3	1:D:113:ILE:CD1	2.24	0.60
1:D:100:ILE:HG12	1:E:103:ILE:HD12	1.84	0.59
1:B:163:TRP:CZ3	1:B:167:VAL:HG11	2.37	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:141:HIS:HE1	1:C:143:TYR:CD2	2.20	0.59
1:D:147:GLU:HG3	1:F:160:TYR:CE2	2.39	0.58
1:D:106:ILE:HG12	1:D:115:GLU:HG2	1.84	0.58
1:A:99:LEU:CD2	1:A:121:MET:HG2	2.34	0.57
1:D:86:ARG:HH12	1:E:113:ILE:HB	1.69	0.57
1:E:99:LEU:CD2	1:E:121:MET:HG2	2.35	0.57
1:E:99:LEU:HD22	1:E:121:MET:HG2	1.87	0.57
1:C:100:ILE:HG22	1:C:103:ILE:HD11	1.86	0.57
1:A:99:LEU:HD22	1:A:121:MET:HG2	1.87	0.56
1:C:158:PHE:HB3	1:C:159:PRO:HD3	1.87	0.56
1:A:162:GLN:HG3	1:A:165:ARG:HH22	1.71	0.56
1:D:88:VAL:HG12	1:D:94:ALA:HB3	1.88	0.54
1:A:135:ASP:OD1	1:A:138:ARG:NH2	2.41	0.54
1:A:165:ARG:O	1:A:169:GLU:HB2	2.07	0.54
1:E:160:TYR:CE2	1:F:147:GLU:HG3	2.42	0.54
1:D:141:HIS:HE1	2:D:207:HOH:O	1.92	0.53
1:A:137:VAL:O	1:A:141:HIS:HB2	2.08	0.53
1:D:109:TRP:CD1	1:D:114:GLU:OE2	2.61	0.53
1:B:141:HIS:HE1	1:B:143:TYR:CD2	2.26	0.53
1:D:147:GLU:HG3	1:F:160:TYR:CZ	2.44	0.52
1:D:162:GLN:O	1:D:166:GLN:HG3	2.09	0.52
1:A:147:GLU:HG3	1:C:160:TYR:CZ	2.45	0.52
1:B:88:VAL:HG12	1:B:94:ALA:HB3	1.92	0.52
1:B:124:LYS:HE3	1:C:147:GLU:OE2	2.10	0.52
1:D:152:PRO:HD3	1:F:154:GLU:HG2	1.91	0.51
1:B:151:LEU:HD12	1:B:151:LEU:N	2.25	0.51
1:D:154:GLU:HG2	1:E:152:PRO:HD3	1.92	0.51
1:A:160:TYR:CE2	1:B:147:GLU:HG3	2.45	0.51
1:C:137:VAL:O	1:C:141:HIS:HB2	2.10	0.51
1:B:97:VAL:HG11	1:B:121:MET:HE3	1.93	0.51
1:F:99:LEU:CD2	1:F:121:MET:HG2	2.40	0.51
1:B:100:ILE:CG2	1:B:103:ILE:HD11	2.38	0.51
1:E:162:GLN:HG3	1:E:165:ARG:HH22	1.75	0.50
1:C:99:LEU:CD2	1:C:121:MET:HG2	2.41	0.50
1:B:98:ASN:ND2	1:B:124:LYS:HZ2	2.10	0.50
1:A:90:GLU:HB2	1:B:113:ILE:HD11	1.93	0.50
1:D:85:ALA:CB	1:E:106:ILE:HD12	2.41	0.50
1:A:87:ALA:O	1:A:91:LYS:HG3	2.12	0.49
1:B:82:LYS:O	1:B:86:ARG:HG3	2.12	0.49
1:D:85:ALA:HB3	1:E:106:ILE:HD12	1.94	0.49
1:E:82:LYS:O	1:E:86:ARG:HG3	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:147:GLU:HG3	1:C:160:TYR:CE2	2.47	0.49
1:B:100:ILE:O	1:B:103:ILE:HD13	2.12	0.49
1:D:157:ASN:OD1	1:D:159:PRO:HD2	2.12	0.49
1:D:109:TRP:CE3	1:D:110:LYS:N	2.80	0.49
1:D:100:ILE:HG12	1:E:103:ILE:CD1	2.42	0.48
1:F:137:VAL:O	1:F:141:HIS:HB2	2.14	0.48
1:D:86:ARG:NH1	1:E:113:ILE:HB	2.26	0.48
1:C:97:VAL:HG11	1:C:121:MET:HE3	1.95	0.48
1:C:162:GLN:HG3	1:C:165:ARG:NH2	2.28	0.48
1:A:109:TRP:CZ2	1:A:110:LYS:HE2	2.48	0.48
1:F:75:CYS:HB3	1:F:80:VAL:HG23	1.95	0.47
2:E:223:HOH:O	1:F:74:THR:HG21	2.13	0.47
1:E:151:LEU:HD23	1:F:151:LEU:HD13	1.96	0.47
1:C:85:ALA:HB2	1:C:121:MET:HE3	1.97	0.47
1:C:87:ALA:O	1:C:91:LYS:HG3	2.14	0.47
1:D:109:TRP:HA	1:D:109:TRP:HE3	1.80	0.47
1:D:150:ALA:C	1:D:151:LEU:HD12	2.35	0.47
1:D:74:THR:HG21	1:D:118:GLU:OE1	2.16	0.46
1:D:109:TRP:CE3	1:D:109:TRP:HA	2.50	0.46
1:A:100:ILE:O	1:A:103:ILE:HD13	2.16	0.46
1:A:100:ILE:HG22	1:A:103:ILE:HD11	1.98	0.46
1:D:151:LEU:HD12	1:D:151:LEU:N	2.31	0.46
1:C:72:PHE:HA	1:C:121:MET:O	2.16	0.45
1:C:111:GLY:C	1:C:112:LYS:HD2	2.36	0.45
1:A:97:VAL:HG11	1:A:121:MET:HE3	1.97	0.45
1:A:151:LEU:HD12	1:A:151:LEU:N	2.31	0.45
1:E:158:PHE:CD1	1:E:158:PHE:N	2.84	0.45
1:F:96:CYS:SG	1:F:124:LYS:HE3	2.57	0.45
1:A:138:ARG:NH1	1:C:157:ASN:OD1	2.50	0.45
1:F:158:PHE:CD1	1:F:158:PHE:N	2.85	0.45
1:A:152:PRO:HD3	1:C:154:GLU:HG2	1.99	0.44
1:A:104:THR:HA	1:A:116:ASP:O	2.17	0.44
1:A:126:GLN:HG3	1:A:168:THR:HG21	2.00	0.44
1:F:99:LEU:HD22	1:F:121:MET:HG2	2.00	0.44
1:F:165:ARG:HH11	1:F:165:ARG:CG	2.31	0.44
1:D:90:GLU:HB2	1:E:113:ILE:HD11	1.99	0.43
1:D:158:PHE:CD1	1:D:158:PHE:N	2.86	0.43
1:D:91:LYS:HE3	1:D:136:PHE:CE1	2.53	0.43
1:B:141:HIS:CE1	1:B:143:TYR:CD2	3.06	0.43
1:F:124:LYS:N	1:F:124:LYS:HD3	2.34	0.43
1:E:73:VAL:HA	1:E:147:GLU:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:75:CYS:HB3	1:F:80:VAL:CG2	2.49	0.43
1:E:97:VAL:HG11	1:E:121:MET:HE3	2.01	0.42
1:A:128:SER:HA	1:D:64:VAL:HG21	2.01	0.42
1:A:154:GLU:HG2	1:B:152:PRO:HD3	2.01	0.42
1:A:157:ASN:HB2	1:B:148:VAL:HB	2.01	0.42
1:A:165:ARG:HG2	1:A:165:ARG:HH11	1.84	0.42
1:D:133:LEU:O	1:D:137:VAL:HG13	2.19	0.42
1:D:151:LEU:N	1:D:151:LEU:CD1	2.82	0.42
1:C:100:ILE:O	1:C:103:ILE:HD13	2.20	0.42
1:D:108:GLU:HG2	1:D:112:LYS:O	2.19	0.42
1:D:88:VAL:CG1	1:D:94:ALA:HB3	2.48	0.42
1:F:98:ASN:HD22	1:F:122:MET:CE	2.33	0.42
1:E:80:VAL:O	1:E:84:ILE:HG13	2.19	0.42
1:D:113:ILE:HG21	1:F:86:ARG:HG2	2.02	0.42
1:B:157:ASN:OD1	1:B:159:PRO:HD2	2.20	0.41
1:E:158:PHE:N	1:E:158:PHE:HD1	2.18	0.41
1:D:99:LEU:CD2	1:D:121:MET:HG2	2.50	0.41
1:B:98:ASN:ND2	1:B:124:LYS:NZ	2.69	0.41
1:B:101:PRO:O	1:B:103:ILE:HD12	2.21	0.41
1:C:124:LYS:N	1:C:124:LYS:HE2	2.36	0.41
1:E:165:ARG:NH1	2:E:191:HOH:O	2.54	0.41
1:F:91:LYS:O	1:F:92:ARG:HB2	2.21	0.41
1:C:86:ARG:HH11	1:C:86:ARG:HB3	1.85	0.40
1:C:129:LEU:CD2	1:C:168:THR:HB	2.51	0.40
1:D:113:ILE:HD11	1:F:90:GLU:CG	2.49	0.40
1:E:75:CYS:O	1:E:118:GLU:HB3	2.22	0.40
1:F:158:PHE:N	1:F:158:PHE:HD1	2.19	0.40
1:C:100:ILE:CG2	1:C:103:ILE:HD11	2.52	0.40
1:A:85:ALA:O	1:A:89:VAL:HG23	2.21	0.40
1:B:88:VAL:CG1	1:B:94:ALA:HB3	2.52	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	107/179 (60%)	104 (97%)	3 (3%)	0	100	100
1	B	105/179 (59%)	105 (100%)	0	0	100	100
1	C	104/179 (58%)	102 (98%)	2 (2%)	0	100	100
1	D	105/179 (59%)	103 (98%)	2 (2%)	0	100	100
1	E	105/179 (59%)	103 (98%)	2 (2%)	0	100	100
1	F	105/179 (59%)	104 (99%)	1 (1%)	0	100	100
All	All	631/1074 (59%)	621 (98%)	10 (2%)	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	95/152 (62%)	93 (98%)	2 (2%)	53	48
1	B	93/152 (61%)	89 (96%)	4 (4%)	29	22
1	C	93/152 (61%)	91 (98%)	2 (2%)	52	46
1	D	93/152 (61%)	89 (96%)	4 (4%)	29	22
1	E	93/152 (61%)	91 (98%)	2 (2%)	52	46
1	F	93/152 (61%)	89 (96%)	4 (4%)	29	22
All	All	560/912 (61%)	542 (97%)	18 (3%)	39	32

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

Mol	Chain	Res	Type
1	A	108	GLU
1	A	124	LYS
1	B	108	GLU
1	B	154	GLU

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Mol	Chain	Res	Type
1	B	161	LEU
1	B	163	TRP
1	C	124	LYS
1	C	129	LEU
1	D	79	LYS
1	D	109	TRP
1	D	124	LYS
1	D	141	HIS
1	E	124	LYS
1	E	154	GLU
1	F	114	GLU
1	F	124	LYS
1	F	129	LEU
1	F	161	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	98	ASN
1	A	162	GLN
1	B	98	ASN
1	B	126	GLN
1	B	141	HIS
1	C	98	ASN
1	C	166	GLN
1	D	98	ASN
1	D	141	HIS
1	E	98	ASN
1	E	102	GLN
1	E	162	GLN
1	F	98	ASN
1	F	126	GLN
1	F	162	GLN

5.3.3 RNA

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	109/179 (60%)	0.28	8 (7%) 15 16	35, 48, 62, 73	0
1	B	107/179 (59%)	0.16	5 (4%) 31 33	34, 47, 64, 77	0
1	C	106/179 (59%)	0.23	6 (5%) 23 25	33, 45, 68, 78	0
1	D	107/179 (59%)	0.68	16 (14%) 2 1	36, 49, 79, 92	0
1	E	107/179 (59%)	0.36	8 (7%) 14 15	33, 49, 78, 98	0
1	F	107/179 (59%)	0.27	4 (3%) 41 45	35, 48, 62, 67	0
All	All	643/1074 (59%)	0.33	47 (7%) 15 16	33, 48, 67, 98	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	109	TRP	7.8
1	D	111	GLY	5.6
1	D	110	LYS	5.4
1	C	108	GLU	4.3
1	E	110	LYS	4.3
1	D	109	TRP	3.9
1	D	112	LYS	3.5
1	A	151	LEU	3.5
1	C	109	TRP	3.4
1	B	151	LEU	3.2
1	F	151	LEU	3.2
1	E	111	GLY	3.2
1	A	61	SER	3.0
1	D	151	LEU	3.0
1	D	114	GLU	2.9
1	C	110	LYS	2.8
1	E	144	GLU	2.8
1	F	109	TRP	2.8
1	A	150	ALA	2.8

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Mol	Chain	Res	Type	RSRZ
1	E	151	LEU	2.8
1	A	149	ILE	2.6
1	D	70	ALA	2.6
1	D	100	ILE	2.6
1	D	152	PRO	2.6
1	D	113	ILE	2.6
1	B	102	GLN	2.5
1	C	153	VAL	2.5
1	B	62	GLY	2.5
1	D	108	GLU	2.5
1	E	108	GLU	2.4
1	D	150	ALA	2.4
1	D	120	LEU	2.4
1	E	145	VAL	2.3
1	B	150	ALA	2.3
1	A	70	ALA	2.3
1	A	158	PHE	2.3
1	A	72	PHE	2.2
1	A	152	PRO	2.2
1	D	62	GLY	2.2
1	B	144	GLU	2.2
1	C	72	PHE	2.2
1	E	153	VAL	2.2
1	F	70	ALA	2.2
1	C	70	ALA	2.1
1	D	71	ALA	2.1
1	F	90	GLU	2.1
1	D	149	ILE	2.0

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.