



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 10:51 PM GMT

PDB ID : 5EGK
Title : The structural and biochemical characterization of acyl-coa hydrolase mutant Asp43Ala from Staphylococcus aureus
Authors : Khandokar, Y.B.; Srivastava, P.S.; Forwood, J.K.
Deposited on : 2015-10-27
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
<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
Mogul : unknown
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
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) : rb-20026982

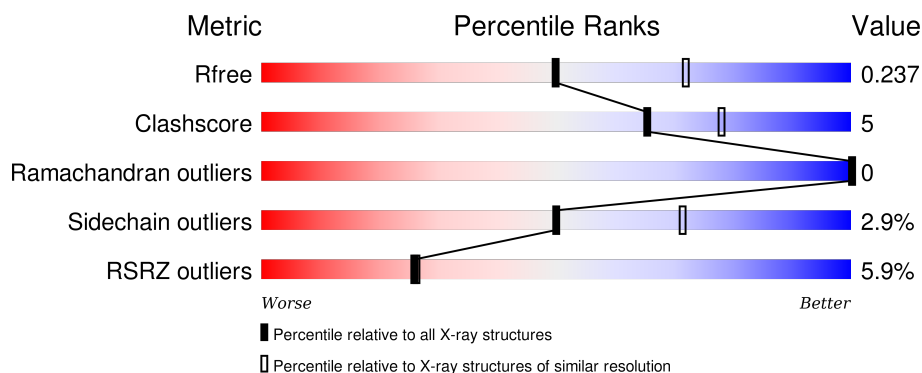
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	91344	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (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 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	179	<div> <div>6%</div> <div>87% 6% 7%</div> </div>
1	B	179	<div> <div>6%</div> <div>77% 13% • 8%</div> </div>
1	C	179	<div> <div>6%</div> <div>84% 7% • 9%</div> </div>
1	D	179	<div> <div>7%</div> <div>83% 9% • 7%</div> </div>
1	E	179	<div> <div>4%</div> <div>78% 13% • 7%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	179	<div><div></div><div>5%</div><div>79%</div><div>12%</div><div>• 8%</div></div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 7831 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 Acyl CoA Hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	166	Total	C	N	O	S	0	0	0
			1311	826	228	249	8			
1	B	164	Total	C	N	O	S	0	0	0
			1295	817	223	247	8			
1	C	163	Total	C	N	O	S	0	0	0
			1292	815	224	245	8			
1	D	167	Total	C	N	O	S	0	0	0
			1319	830	229	252	8			
1	E	167	Total	C	N	O	S	0	0	0
			1319	830	229	252	8			
1	F	164	Total	C	N	O	S	0	0	0
			1295	817	223	247	8			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP A0A0H3K033
A	-1	ASN	-	expression tag	UNP A0A0H3K033
A	0	ALA	-	expression tag	UNP A0A0H3K033
A	43	ALA	ASP	engineered mutation	UNP A0A0H3K033
B	-2	SER	-	expression tag	UNP A0A0H3K033
B	-1	ASN	-	expression tag	UNP A0A0H3K033
B	0	ALA	-	expression tag	UNP A0A0H3K033
B	43	ALA	ASP	engineered mutation	UNP A0A0H3K033
C	-2	SER	-	expression tag	UNP A0A0H3K033
C	-1	ASN	-	expression tag	UNP A0A0H3K033
C	0	ALA	-	expression tag	UNP A0A0H3K033
C	43	ALA	ASP	engineered mutation	UNP A0A0H3K033
D	-2	SER	-	expression tag	UNP A0A0H3K033
D	-1	ASN	-	expression tag	UNP A0A0H3K033
D	0	ALA	-	expression tag	UNP A0A0H3K033
D	43	ALA	ASP	engineered mutation	UNP A0A0H3K033
E	-2	SER	-	expression tag	UNP A0A0H3K033

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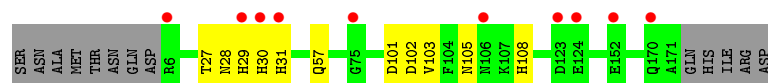
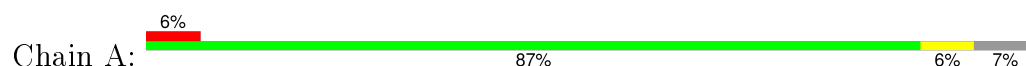
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Chain	Residue	Modelled	Actual	Comment	Reference
E	-1	ASN	-	expression tag	UNP A0A0H3K033
E	0	ALA	-	expression tag	UNP A0A0H3K033
E	43	ALA	ASP	engineered mutation	UNP A0A0H3K033
F	-2	SER	-	expression tag	UNP A0A0H3K033
F	-1	ASN	-	expression tag	UNP A0A0H3K033
F	0	ALA	-	expression tag	UNP A0A0H3K033
F	43	ALA	ASP	engineered mutation	UNP A0A0H3K033

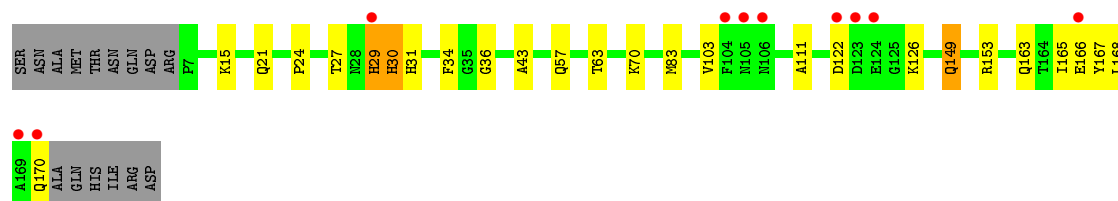
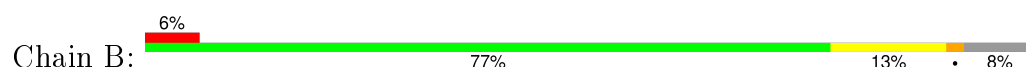
3 Residue-property plots [i](#)

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.

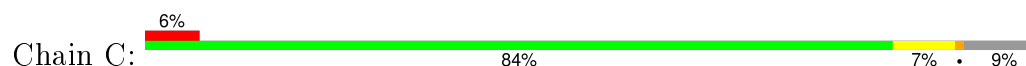
• Molecule 1: Acyl CoA Hydrolase



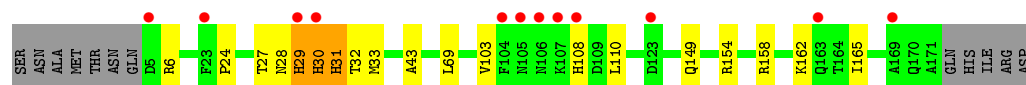
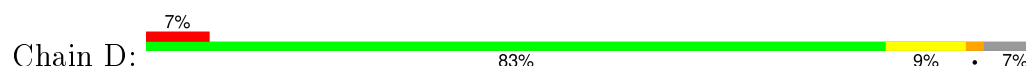
• Molecule 1: Acyl CoA Hydrolase



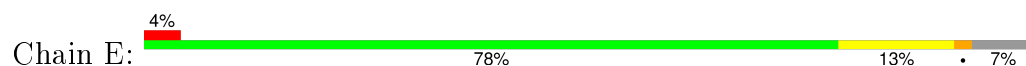
• Molecule 1: Acyl CoA Hydrolase



• Molecule 1: Acyl CoA Hydrolase

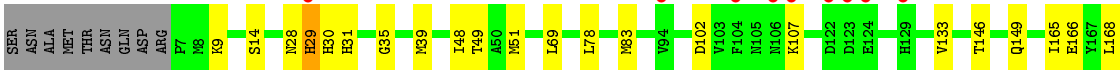
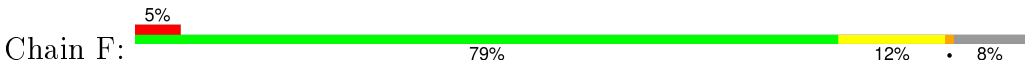


• Molecule 1: Acyl CoA Hydrolase



ILE
ARG
ASP

● Molecule 1: Acyl CoA Hydrolase



A169
Q170
ALA
GLN
HIS
ILE
ARG
ASP

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	49.71 Å 129.00 Å 83.56 Å 90.00° 104.10° 90.00°	Depositor
Resolution (Å)	29.96 – 2.40 29.96 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.8 (29.96-2.40) 98.9 (29.96-2.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.94 (at 2.39 Å)	Xtriage
Refinement program	PHENIX (1.10pre_2104: ???)	Depositor
R, R_{free}	0.206 , 0.238 0.205 , 0.237	Depositor DCC
R_{free} test set	1920 reflections (4.87%)	DCC
Wilson B-factor (Å ²)	37.3	Xtriage
Anisotropy	0.047	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 39.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	3 of 39444 reflections (0.008%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7831	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 4.24% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.37	0/1338	0.54	0/1810
1	B	0.40	0/1322	0.56	0/1788
1	C	0.40	0/1319	0.56	0/1784
1	D	0.39	0/1346	0.58	0/1821
1	E	0.40	0/1346	0.56	0/1821
1	F	0.41	1/1322 (0.1%)	0.53	0/1788
All	All	0.39	1/7993 (0.0%)	0.56	0/10812

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	35	GLY	C-N	-6.34	1.21	1.33

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	1311	0	1296	6	0
1	B	1295	0	1279	18	0
1	C	1292	0	1278	12	0
1	D	1319	0	1300	22	0
1	E	1319	0	1300	25	0
1	F	1295	0	1279	14	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	7831	0	7732	78	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 5.

All (78) 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:29:HIS:CD2	1:D:30:HIS:CE1	2.48	1.02
1:D:29:HIS:HD2	1:D:30:HIS:ND1	1.57	1.00
1:D:29:HIS:CD2	1:D:30:HIS:ND1	2.36	0.92
1:E:28:ASN:OD1	1:E:30:HIS:N	2.01	0.92
1:F:28:ASN:OD1	1:F:30:HIS:N	2.09	0.85
1:E:24:PRO:O	1:E:27:THR:OG1	1.95	0.84
1:E:28:ASN:HD21	1:E:32:THR:HB	1.56	0.70
1:A:28:ASN:OD1	1:A:30:HIS:N	2.27	0.68
1:B:30:HIS:O	1:B:31:HIS:HB2	1.95	0.65
1:F:30:HIS:O	1:F:31:HIS:HB2	1.97	0.65
1:D:29:HIS:HD2	1:D:30:HIS:CE1	1.99	0.65
1:C:30:HIS:O	1:C:31:HIS:HB2	1.96	0.64
1:D:69:LEU:HD13	1:F:165:ILE:HD11	1.80	0.64
1:B:24:PRO:O	1:B:27:THR:OG1	2.17	0.63
1:D:30:HIS:O	1:D:31:HIS:HB2	2.00	0.61
1:C:146:THR:O	1:C:149:GLN:HG2	2.01	0.60
1:B:167:TYR:O	1:B:170:GLN:HG2	2.03	0.59
1:D:110:LEU:HD22	1:F:168:LEU:HD12	1.86	0.58
1:E:148:PRO:O	1:E:152:GLU:HG3	2.05	0.57
1:C:21:GLN:NE2	1:F:29:HIS:CE1	2.72	0.57
1:B:21:GLN:NE2	1:D:29:HIS:CE1	2.74	0.55
1:B:21:GLN:NE2	1:D:29:HIS:HE1	2.03	0.55
1:C:21:GLN:NE2	1:F:29:HIS:NE2	2.57	0.53
1:B:57:GLN:HA	1:B:57:GLN:OE1	2.09	0.52
1:E:30:HIS:O	1:E:31:HIS:HB2	2.10	0.51
1:B:70:LYS:HB2	1:B:111:ALA:HA	1.93	0.51
1:E:28:ASN:OD1	1:E:30:HIS:HB2	2.11	0.50
1:D:24:PRO:O	1:D:27:THR:OG1	2.28	0.50
1:A:102:ASP:OD2	1:A:105:ASN:ND2	2.45	0.49
1:C:29:HIS:ND1	1:C:29:HIS:N	2.60	0.49
1:C:29:HIS:HB3	1:E:74:THR:HG22	1.94	0.48
1:B:165:ILE:HG12	1:E:110:LEU:HD23	1.95	0.48
1:D:30:HIS:N	1:D:30:HIS:ND1	2.60	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:149:GLN:HA	1:E:152:GLU:OE1	2.13	0.48
1:D:165:ILE:HD11	1:F:69:LEU:HD22	1.95	0.48
1:B:15:LYS:HE2	1:B:83:MET:CE	2.43	0.48
1:D:28:ASN:HD21	1:D:32:THR:HB	1.79	0.48
1:C:16:CYS:SG	1:C:52:LYS:HD2	2.54	0.48
1:C:28:ASN:OD1	1:C:30:HIS:N	2.47	0.47
1:F:14:SER:O	1:F:49:THR:HG23	2.15	0.47
1:A:57:GLN:NE2	1:C:30:HIS:HE1	2.13	0.47
1:E:70:LYS:HB2	1:E:111:ALA:HA	1.97	0.47
1:C:55:GLY:HA2	1:D:103:VAL:HG12	1.97	0.47
1:F:146:THR:O	1:F:149:GLN:HG2	2.16	0.46
1:F:102:ASP:HB3	1:F:107:LYS:HB3	1.97	0.46
1:F:28:ASN:OD1	1:F:30:HIS:HB2	2.17	0.45
1:D:162:LYS:O	1:D:165:ILE:HB	2.17	0.45
1:B:122:ASP:OD2	1:B:126:LYS:HB2	2.17	0.45
1:E:101:ASP:OD1	1:E:108:HIS:CD2	2.70	0.45
1:E:101:ASP:OD1	1:E:108:HIS:NE2	2.50	0.45
1:E:122:ASP:HB3	1:E:128:LYS:HD2	1.99	0.45
1:E:101:ASP:OD1	1:E:108:HIS:CE1	2.71	0.44
1:B:34:PHE:CE2	1:B:36:GLY:HA3	2.53	0.44
1:E:157:ARG:HA	1:E:160:GLU:HG2	1.99	0.44
1:C:70:LYS:HE3	1:C:70:LYS:HB3	1.79	0.43
1:B:43:ALA:HB2	1:E:39:MET:HE1	2.00	0.43
1:B:29:HIS:N	1:B:29:HIS:ND1	2.60	0.43
1:D:6:ARG:HA	1:D:6:ARG:HD3	1.67	0.43
1:B:63:THR:HB	1:E:66:VAL:HB	2.00	0.42
1:D:32:THR:HG22	1:D:33:MET:N	2.33	0.42
1:F:9:LYS:O	1:F:133:VAL:HA	2.19	0.42
1:E:102:ASP:HB3	1:E:105:ASN:OD1	2.18	0.42
1:B:168:LEU:HD12	1:E:110:LEU:HD22	2.02	0.42
1:A:101:ASP:OD2	1:E:52:LYS:NZ	2.34	0.42
1:E:28:ASN:ND2	1:E:32:THR:HB	2.30	0.42
1:C:52:LYS:HA	1:D:103:VAL:HG13	2.01	0.42
1:B:149:GLN:OE1	1:B:153:ARG:NH1	2.52	0.42
1:D:162:LYS:HA	1:D:165:ILE:HD12	2.02	0.41
1:B:163:GLN:O	1:B:166:GLU:HG2	2.20	0.41
1:F:48:ILE:HA	1:F:51:MET:HE2	2.02	0.41
1:E:100:ILE:HG13	1:E:111:ALA:HB2	2.02	0.41
1:E:70:LYS:HE3	1:E:70:LYS:HB3	1.77	0.41
1:A:30:HIS:O	1:A:31:HIS:HB2	2.20	0.41
1:E:30:HIS:O	1:E:31:HIS:CB	2.69	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:43:ALA:HB2	1:F:39:MET:CE	2.51	0.40
1:A:27:THR:O	1:D:24:PRO:HD2	2.21	0.40
1:D:154:ARG:O	1:D:158:ARG:HG3	2.21	0.40
1:B:43:ALA:HB2	1:E:39:MET:CE	2.50	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	164/179 (92%)	163 (99%)	1 (1%)	0	100	100
1	B	162/179 (90%)	162 (100%)	0	0	100	100
1	C	161/179 (90%)	160 (99%)	1 (1%)	0	100	100
1	D	165/179 (92%)	164 (99%)	1 (1%)	0	100	100
1	E	165/179 (92%)	164 (99%)	1 (1%)	0	100	100
1	F	162/179 (90%)	161 (99%)	1 (1%)	0	100	100
All	All	979/1074 (91%)	974 (100%)	5 (0%)	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	142/154 (92%)	139 (98%)	3 (2%)	61	80
1	B	141/154 (92%)	137 (97%)	4 (3%)	51	72
1	C	141/154 (92%)	138 (98%)	3 (2%)	61	80
1	D	143/154 (93%)	138 (96%)	5 (4%)	43	64
1	E	143/154 (93%)	137 (96%)	6 (4%)	36	56
1	F	141/154 (92%)	137 (97%)	4 (3%)	51	72
All	All	851/924 (92%)	826 (97%)	25 (3%)	50	71

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	HIS
1	A	103	VAL
1	A	108	HIS
1	B	29	HIS
1	B	30	HIS
1	B	103	VAL
1	B	149	GLN
1	C	6	ARG
1	C	31	HIS
1	C	103	VAL
1	D	29	HIS
1	D	30	HIS
1	D	31	HIS
1	D	108	HIS
1	D	149	GLN
1	E	29	HIS
1	E	30	HIS
1	E	31	HIS
1	E	83	MET
1	E	101	ASP
1	E	108	HIS
1	F	29	HIS
1	F	78	LEU
1	F	83	MET
1	F	166	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	HIS
1	A	57	GLN
1	B	21	GLN
1	C	21	GLN
1	C	30	HIS
1	D	29	HIS
1	D	31	HIS
1	E	31	HIS
1	E	57	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 carbohydrates 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 ⓘ

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	166/179 (92%)	0.37	10 (6%)	25 25	20, 36, 67, 88	0
1	B	164/179 (91%)	0.30	10 (6%)	25 25	25, 40, 73, 88	0
1	C	163/179 (91%)	0.42	10 (6%)	25 25	20, 36, 66, 88	0
1	D	167/179 (93%)	0.40	12 (7%)	18 18	23, 35, 73, 88	0
1	E	167/179 (93%)	0.34	7 (4%)	40 41	22, 37, 70, 88	0
1	F	164/179 (91%)	0.41	9 (5%)	29 29	25, 42, 74, 88	0
All	All	991/1074 (92%)	0.37	58 (5%)	26 26	20, 38, 71, 88	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	29	HIS	6.4
1	C	29	HIS	5.6
1	B	29	HIS	5.4
1	D	29	HIS	5.2
1	A	29	HIS	5.0
1	B	124	GLU	4.7
1	D	105	ASN	4.7
1	D	30	HIS	4.2
1	A	31	HIS	4.1
1	B	104	PHE	4.0
1	F	29	HIS	3.6
1	F	123	ASP	3.6
1	E	159	ILE	3.4
1	B	123	ASP	3.4
1	F	104	PHE	3.3
1	C	159	ILE	3.3
1	E	152	GLU	3.3
1	E	30	HIS	3.1
1	E	107	LYS	3.1

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Mol	Chain	Res	Type	RSRZ
1	D	23	PHE	3.1
1	A	152	GLU	3.0
1	D	123	ASP	3.0
1	F	107	LYS	3.0
1	D	169	ALA	2.9
1	F	106	ASN	2.9
1	A	123	ASP	2.8
1	D	5	ASP	2.8
1	A	106	ASN	2.8
1	C	123	ASP	2.7
1	C	106	ASN	2.7
1	B	170	GLN	2.7
1	F	124	GLU	2.7
1	C	30	HIS	2.7
1	D	108	HIS	2.6
1	A	6	ARG	2.6
1	A	30	HIS	2.6
1	C	138	ASP	2.6
1	B	122	ASP	2.5
1	D	107	LYS	2.5
1	D	104	PHE	2.5
1	B	105	ASN	2.4
1	D	106	ASN	2.4
1	C	125	GLY	2.4
1	C	155	LYS	2.3
1	B	106	ASN	2.3
1	B	169	ALA	2.3
1	A	75	GLY	2.2
1	E	104	PHE	2.2
1	E	108	HIS	2.2
1	A	170	GLN	2.2
1	A	124	GLU	2.2
1	C	107	LYS	2.1
1	F	122	ASP	2.1
1	F	129	HIS	2.1
1	C	153	ARG	2.0
1	F	94	VAL	2.0
1	B	166	GLU	2.0
1	D	163	GLN	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 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.