



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:08 AM GMT

PDB ID : 3DIF
Title : Crystal structure of FabOX117
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Deposited on : 2008-06-20
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 : 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

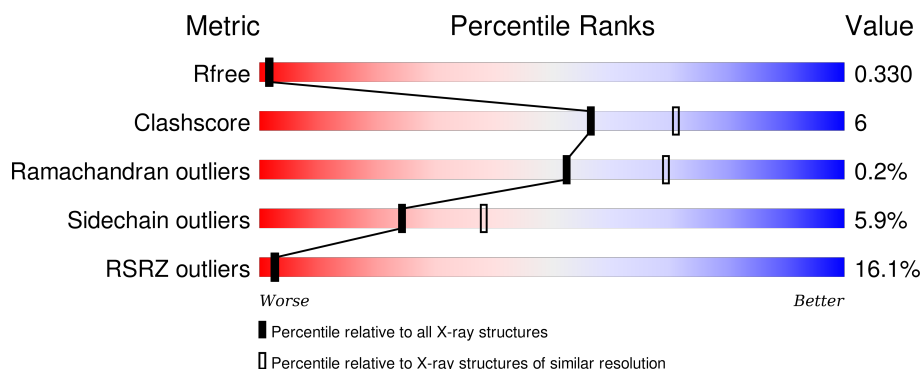
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	214	<div> <div>17%</div> <div>82%</div> <div>16%</div> <div>.</div> </div>
1	C	214	<div> <div>25%</div> <div>83%</div> <div>16%</div> <div>.</div> </div>
2	B	229	<div> <div>9%</div> <div>75%</div> <div>14%</div> <div>.</div> <div>8%</div> </div>
2	D	229	<div> <div>12%</div> <div>82%</div> <div>9%</div> <div>.</div> <div>7%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6854 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 FabOX117 Light Chain Fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	213	Total	C	N	O	S	0	0	0
			1657	1038	276	337	6			
1	C	214	Total	C	N	O	S	0	0	0
			1663	1041	277	338	7			

- Molecule 2 is a protein called FabOX117 Heavy Chain Fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	211	Total	C	N	O	S	0	0	0
			1604	1030	259	309	6			
2	D	213	Total	C	N	O	S	0	0	0
			1619	1038	264	311	6			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	224	HIS	-	EXPRESSION TAG	PDB 3DIF
B	225	HIS	-	EXPRESSION TAG	PDB 3DIF
B	226	HIS	-	EXPRESSION TAG	PDB 3DIF
B	227	HIS	-	EXPRESSION TAG	PDB 3DIF
B	228	HIS	-	EXPRESSION TAG	PDB 3DIF
B	229	HIS	-	EXPRESSION TAG	PDB 3DIF
D	224	HIS	-	EXPRESSION TAG	PDB 3DIF
D	225	HIS	-	EXPRESSION TAG	PDB 3DIF
D	226	HIS	-	EXPRESSION TAG	PDB 3DIF
D	227	HIS	-	EXPRESSION TAG	PDB 3DIF
D	228	HIS	-	EXPRESSION TAG	PDB 3DIF
D	229	HIS	-	EXPRESSION TAG	PDB 3DIF

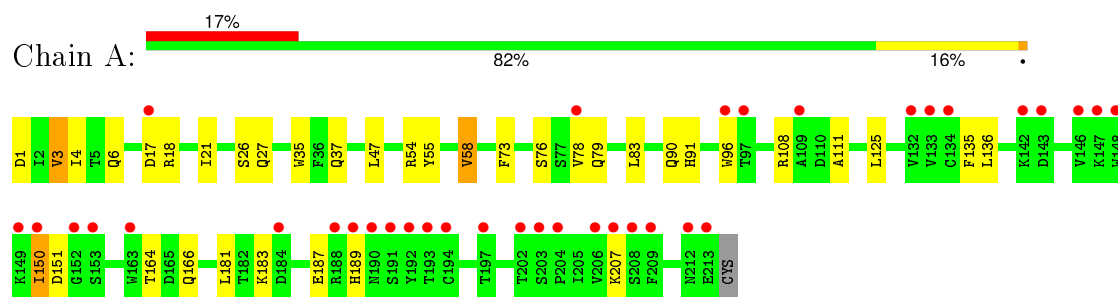
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	39	Total 39	O 39	0	0
3	B	83	Total 83	O 83	0	0
3	C	74	Total 74	O 74	0	0
3	D	115	Total 115	O 115	0	0

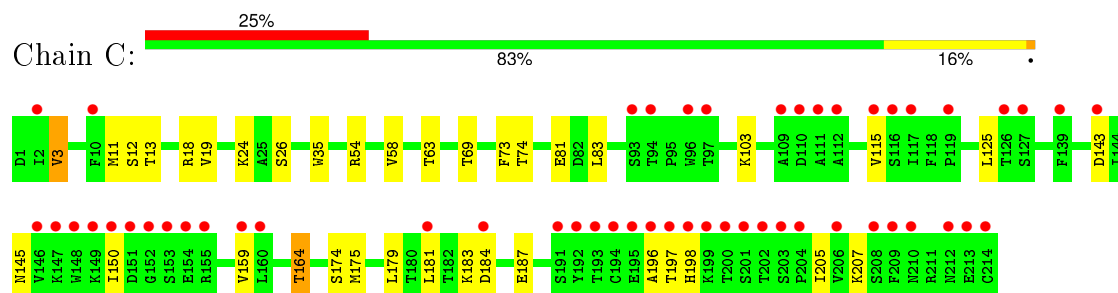
3 Residue-property plots [i](#)

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.

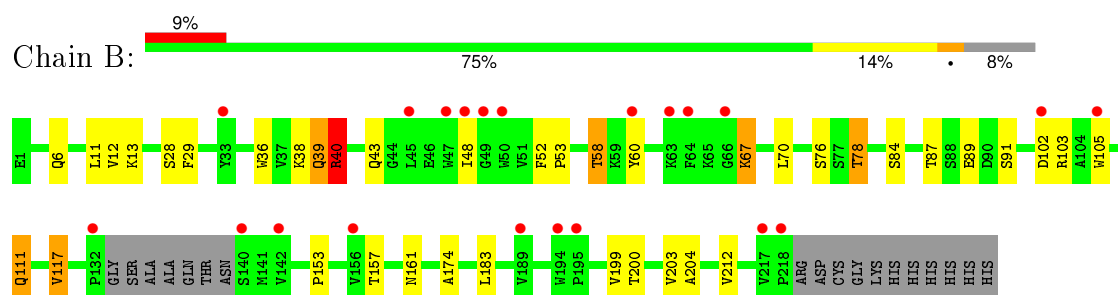
- Molecule 1: FabOX117 Light Chain Fragment



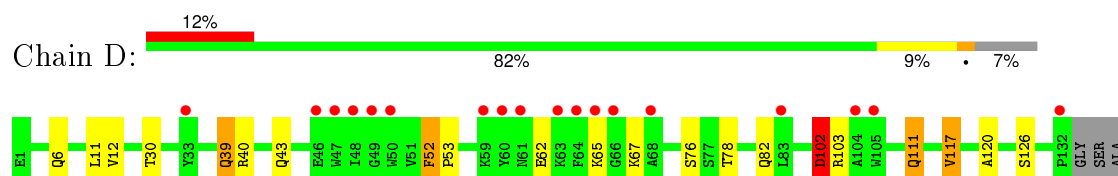
- Molecule 1: FabOX117 Light Chain Fragment

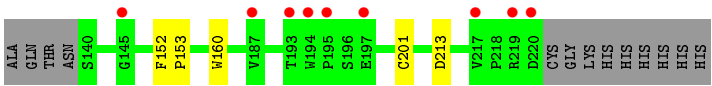


- Molecule 2: FabOX117 Heavy Chain Fragment



- Molecule 2: FabOX117 Heavy Chain Fragment





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	41.21Å 70.60Å 87.85Å 101.50° 102.94° 98.32°	Depositor
Resolution (Å)	30.00 – 2.40 33.91 – 2.37	Depositor EDS
% Data completeness (in resolution range)	97.7 (30.00-2.40) 90.0 (33.91-2.37)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.63 (at 2.36Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.219 , 0.286 0.284 , 0.330	Depositor DCC
R_{free} test set	1777 reflections (5.02%)	DCC
Wilson B-factor (Å ²)	34.7	Xtriage
Anisotropy	0.354	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 38.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 35858 reflections	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	6854	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.33% 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	0.35	0/1697	0.55	0/2307
1	C	0.36	0/1703	0.53	0/2315
2	B	0.41	0/1652	0.58	1/2259 (0.0%)
2	D	0.38	0/1667	0.54	0/2278
All	All	0.38	0/6719	0.55	1/9159 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	40	ARG	NE-CZ-NH1	6.79	123.69	120.30

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	1657	0	1583	22	0
1	C	1663	0	1588	16	0
2	B	1604	0	1569	27	0
2	D	1619	0	1583	17	0
3	A	39	0	0	0	0
3	B	83	0	0	3	0
3	C	74	0	0	0	0
3	D	115	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	6854	0	6323	79	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 6.

All (79) 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)
2:D:111:GLN:HE21	2:D:111:GLN:H	1.19	0.86
2:B:87:THR:HG22	2:B:89:GLU:H	1.39	0.86
2:B:6:GLN:H	2:B:111:GLN:HE22	1.19	0.84
2:D:6:GLN:H	2:D:111:GLN:HE22	1.25	0.84
2:B:111:GLN:H	2:B:111:GLN:HE21	1.34	0.74
1:A:54:ARG:HG3	1:A:58:VAL:HG22	1.68	0.73
2:D:39:GLN:HG2	3:D:328:HOH:O	1.90	0.69
1:A:3:VAL:HG13	1:A:26:SER:HB3	1.74	0.68
2:B:40:ARG:HH11	2:B:40:ARG:HG3	1.60	0.67
2:B:203:VAL:HB	2:B:212:VAL:HG23	1.76	0.66
3:B:277:HOH:O	2:D:102:ASP:HB3	1.97	0.66
1:A:17:ASP:O	1:A:78:VAL:HG23	1.96	0.66
2:D:102:ASP:OD1	3:D:238:HOH:O	2.12	0.65
1:A:83:LEU:HD11	1:A:166:GLN:HB3	1.78	0.65
1:A:27:GLN:OE1	2:D:30:THR:HG21	1.97	0.63
1:C:54:ARG:HH12	1:C:63:THR:HG22	1.63	0.63
1:C:13:THR:HG21	1:C:19:VAL:HG22	1.80	0.62
2:D:76:SER:OG	2:D:78:THR:HG22	2.00	0.60
1:A:37:GLN:HB2	1:A:47:LEU:HD11	1.84	0.60
2:B:76:SER:O	2:B:78:THR:HG22	2.03	0.59
2:B:40:ARG:HB2	2:B:43:GLN:HB2	1.85	0.58
1:A:150:ILE:HD11	1:A:189:HIS:CG	2.39	0.58
1:C:3:VAL:HG13	1:C:26:SER:HB3	1.85	0.57
2:D:12:VAL:O	2:D:117:VAL:HA	2.05	0.57
1:C:150:ILE:HD11	1:C:179:LEU:HD21	1.85	0.56
1:C:196:ALA:HB3	1:C:205:ILE:HG23	1.87	0.56
1:A:54:ARG:CG	1:A:58:VAL:HG22	2.34	0.55
1:C:145:ASN:HB2	1:C:197:THR:HB	1.89	0.55
1:C:196:ALA:HB3	1:C:205:ILE:CG2	2.38	0.54
1:A:6:GLN:HE21	1:A:21:ILE:HG21	1.73	0.54
2:B:87:THR:O	2:B:117:VAL:HG11	2.08	0.54
2:B:76:SER:OG	2:B:78:THR:HG23	2.08	0.53
2:D:111:GLN:N	2:D:111:GLN:HE21	1.97	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:18:ARG:HG3	1:A:76:SER:HA	1.91	0.53
1:A:4:ILE:HD11	1:A:90:GLN:HB3	1.92	0.52
2:B:38:LYS:HB2	2:B:48:ILE:HD11	1.91	0.51
2:B:12:VAL:O	2:B:117:VAL:HA	2.12	0.50
2:B:6:GLN:H	2:B:111:GLN:NE2	1.99	0.50
1:C:159:VAL:HG22	1:C:179:LEU:HD13	1.93	0.49
1:C:54:ARG:NH1	1:C:63:THR:HG22	2.26	0.49
2:B:103:ARG:HD3	2:B:105:TRP:CH2	2.47	0.49
2:B:40:ARG:HH11	2:B:40:ARG:CG	2.25	0.49
2:B:11:LEU:HB2	2:B:153:PRO:HG3	1.94	0.49
2:B:161:ASN:HD21	2:B:199:VAL:HA	1.78	0.49
1:C:115:VAL:HG23	1:C:205:ILE:HD13	1.95	0.48
2:B:161:ASN:ND2	2:B:200:THR:H	2.11	0.48
2:B:39:GLN:HG2	3:B:285:HOH:O	2.14	0.48
1:A:55:TYR:O	1:A:58:VAL:HG13	2.13	0.48
2:D:40:ARG:HB2	2:D:43:GLN:HB2	1.94	0.48
1:A:1:ASP:N	2:D:53:PRO:O	2.46	0.48
1:C:164:THR:HG22	1:C:174:SER:H	1.79	0.47
2:D:111:GLN:H	2:D:111:GLN:NE2	2.00	0.47
1:A:35:TRP:CE2	1:A:73:PHE:HB2	2.49	0.47
2:B:103:ARG:HD3	2:B:105:TRP:CZ3	2.49	0.47
2:D:11:LEU:HB2	2:D:153:PRO:HG3	1.96	0.47
1:A:1:ASP:HB2	2:D:52:PHE:CE2	2.50	0.46
1:C:35:TRP:CE2	1:C:73:PHE:HB2	2.51	0.46
2:B:157:THR:CG2	2:B:204:ALA:HB3	2.46	0.45
1:C:11:MET:CE	1:C:12:SER:H	2.29	0.45
1:A:54:ARG:HG3	1:A:58:VAL:CG2	2.43	0.45
1:A:136:LEU:HD12	1:A:136:LEU:N	2.32	0.44
1:A:183:LYS:O	1:A:187:GLU:HG2	2.17	0.44
2:D:62:GLU:HA	2:D:65:LYS:HD3	2.01	0.43
2:D:160:TRP:CZ3	2:D:201:CYS:HB3	2.53	0.43
2:B:103:ARG:HB2	3:B:296:HOH:O	2.19	0.43
2:B:29:PHE:CE2	2:B:53:PRO:HB3	2.53	0.43
1:C:13:THR:HG21	1:C:19:VAL:CG2	2.46	0.43
2:B:91:SER:OG	2:B:117:VAL:HG13	2.19	0.42
1:C:183:LYS:O	1:C:187:GLU:HG2	2.20	0.42
2:B:174:ALA:HB2	2:B:183:LEU:HD23	2.02	0.41
1:A:108:ARG:HH21	1:A:111:ALA:HB2	1.85	0.41
2:B:58:THR:HG23	2:B:60:TYR:CE2	2.55	0.41
1:A:91:HIS:HA	1:A:96:TRP:CD1	2.56	0.41
1:A:135:PHE:C	1:A:136:LEU:HD12	2.41	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:18:ARG:CZ	1:C:74:THR:HG21	2.52	0.40
2:B:36:TRP:CD1	2:B:70:LEU:HD22	2.57	0.40
1:A:150:ILE:HD11	1:A:189:HIS:HB3	2.04	0.40
2:B:67:LYS:HE2	2:B:84:SER:O	2.22	0.40
2:D:120:ALA:HB3	2:D:152:PHE:CE2	2.57	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	211/214 (99%)	205 (97%)	6 (3%)	0	100	100
1	C	212/214 (99%)	205 (97%)	7 (3%)	0	100	100
2	B	207/229 (90%)	202 (98%)	4 (2%)	1 (0%)	34	48
2	D	209/229 (91%)	205 (98%)	3 (1%)	1 (0%)	34	48
All	All	839/886 (95%)	817 (97%)	20 (2%)	2 (0%)	52	69

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	102	ASP
2	B	102	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	189/190 (100%)	180 (95%)	9 (5%)	31	49
1	C	190/190 (100%)	175 (92%)	15 (8%)	15	23
2	B	180/194 (93%)	170 (94%)	10 (6%)	26	41
2	D	181/194 (93%)	171 (94%)	10 (6%)	27	42
All	All	740/768 (96%)	696 (94%)	44 (6%)	24	38

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	VAL
1	A	58	VAL
1	A	79	GLN
1	A	125	LEU
1	A	150	ILE
1	A	151	ASP
1	A	164	THR
1	A	181	LEU
1	A	207	LYS
2	B	13	LYS
2	B	28	SER
2	B	39	GLN
2	B	40	ARG
2	B	52	PHE
2	B	58	THR
2	B	67	LYS
2	B	78	THR
2	B	111	GLN
2	B	117	VAL
1	C	3	VAL
1	C	24	LYS
1	C	58	VAL
1	C	69	THR
1	C	81	GLU
1	C	83	LEU
1	C	103	LYS
1	C	125	LEU
1	C	143	ASP
1	C	164	THR
1	C	175	MET
1	C	181	LEU

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Mol	Chain	Res	Type
1	C	184	ASP
1	C	198	HIS
1	C	207	LYS
2	D	39	GLN
2	D	52	PHE
2	D	67	LYS
2	D	82	GLN
2	D	102	ASP
2	D	103	ARG
2	D	111	GLN
2	D	117	VAL
2	D	126	SER
2	D	213	ASP

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	145	ASN
2	B	111	GLN
2	B	161	ASN
1	C	137	ASN
1	C	161	ASN
1	C	198	HIS
2	D	5	GLN
2	D	82	GLN
2	D	111	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	213/214 (99%)	1.08	36 (16%) 2 2	40, 53, 64, 70	0
1	C	214/214 (100%)	1.19	53 (24%) 1 1	36, 54, 65, 82	0
2	B	211/229 (92%)	0.77	21 (9%) 9 9	42, 54, 67, 82	0
2	D	213/229 (93%)	0.79	27 (12%) 5 5	41, 54, 69, 80	0
All	All	851/886 (96%)	0.96	137 (16%) 3 2	36, 54, 67, 82	0

All (137) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	214	CYS	7.6
1	C	200	THR	7.0
2	B	194	TRP	6.3
1	C	206	VAL	5.9
1	C	201	SER	5.3
1	C	193	THR	5.2
1	C	153	SER	5.0
1	C	202	THR	4.9
1	A	197	THR	4.7
1	A	194	CYS	4.7
1	C	194	CYS	4.6
1	A	150	ILE	4.5
1	C	184	ASP	4.4
1	C	148	TRP	4.3
1	C	197	THR	4.2
1	A	152	GLY	4.1
1	A	206	VAL	4.1
1	A	134	CYS	4.1
1	A	213	GLU	3.9
1	C	203	SER	3.9
1	C	198	HIS	3.8

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Mol	Chain	Res	Type	RSRZ
1	C	147	LYS	3.8
1	A	191	SER	3.8
1	A	203	SER	3.8
1	A	190	ASN	3.8
2	D	194	TRP	3.7
1	A	204	PRO	3.7
1	A	153	SER	3.7
2	D	49	GLY	3.6
2	D	197	GLU	3.6
2	D	48	ILE	3.6
1	C	149	LYS	3.5
2	D	193	THR	3.5
1	A	208	SER	3.5
1	A	133	VAL	3.5
2	D	47	TRP	3.4
2	D	132	PRO	3.4
1	C	195	GLU	3.4
2	D	220	ASP	3.4
1	C	212	ASN	3.4
1	A	147	LYS	3.3
1	A	212	ASN	3.3
2	D	64	PHE	3.3
2	D	219	ARG	3.3
1	C	115	VAL	3.3
2	D	61	ASN	3.3
1	A	202	THR	3.2
1	C	192	TYR	3.2
1	C	209	PHE	3.2
2	B	47	TRP	3.2
1	C	127	SER	3.2
1	C	117	ILE	3.2
1	C	208	SER	3.1
1	C	213	GLU	3.1
1	C	210	ASN	3.1
2	B	132	PRO	3.0
2	D	60	TYR	3.0
2	D	195	PRO	3.0
1	C	191	SER	2.8
1	A	17	ASP	2.8
1	A	193	THR	2.7
2	B	64	PHE	2.7
1	A	78	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	146	VAL	2.7
1	C	155	ARG	2.7
1	C	159	VAL	2.6
1	A	146	VAL	2.6
1	A	143	ASP	2.6
1	A	207	LYS	2.6
2	B	60	TYR	2.6
1	A	163	TRP	2.5
2	D	104	ALA	2.5
1	C	139	PHE	2.5
2	B	142	VAL	2.5
1	A	148	TRP	2.5
1	A	184	ASP	2.5
1	C	112	ALA	2.5
2	B	49	GLY	2.5
1	C	111	ALA	2.5
1	C	10	PHE	2.5
1	A	188	ARG	2.5
2	B	189	VAL	2.5
1	C	160	LEU	2.5
1	A	149	LYS	2.5
1	A	132	VAL	2.5
2	B	48	ILE	2.5
1	C	96	TRP	2.4
2	B	50	TRP	2.4
2	D	50	TRP	2.4
2	D	66	GLY	2.4
2	B	195	PRO	2.4
2	D	217	VAL	2.4
2	B	33	TYR	2.4
1	C	97	THR	2.3
2	D	63	LYS	2.3
1	C	154	GLU	2.3
1	A	109	ALA	2.3
2	D	65	LYS	2.3
1	C	116	SER	2.3
2	B	140	SER	2.3
1	A	189	HIS	2.3
1	C	151	ASP	2.3
1	C	109	ALA	2.3
1	A	209	PHE	2.3
1	A	96	TRP	2.2

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Mol	Chain	Res	Type	RSRZ
2	B	66	GLY	2.2
1	A	192	TYR	2.2
2	B	102	ASP	2.2
1	C	199	LYS	2.2
2	D	105	TRP	2.2
1	C	2	ILE	2.2
1	C	126	THR	2.2
2	B	217	VAL	2.2
1	A	97	THR	2.2
2	D	68	ALA	2.1
1	C	143	ASP	2.1
2	B	45	LEU	2.1
1	C	94	THR	2.1
1	C	93	SER	2.1
1	C	196	ALA	2.1
2	B	156	VAL	2.1
1	C	181	LEU	2.1
1	A	142	LYS	2.1
2	D	33	TYR	2.1
2	D	83	LEU	2.1
2	D	187	VAL	2.1
2	B	63	LYS	2.1
2	D	59	LYS	2.0
1	C	119	PRO	2.0
2	B	218	PRO	2.0
1	C	152	GLY	2.0
2	D	145	GLY	2.0
1	C	110	ASP	2.0
1	C	150	ILE	2.0
1	C	204	PRO	2.0
2	D	46	GLU	2.0
2	B	105	TRP	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.