



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 03:04 am GMT

PDB ID : 4ELO
Title : Ferric binding protein in apo form 1
Authors : Wang, Q.; Liu, X.Q.; Wang, X.Q.
Deposited on : 2012-04-11
Resolution : 1.91 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
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

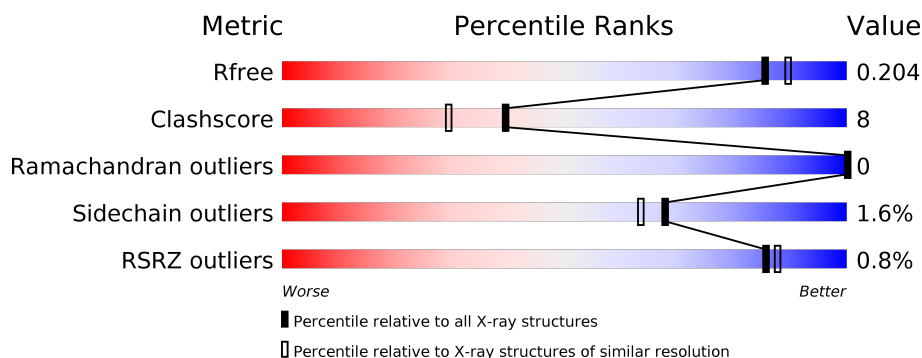
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.91 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}	100719	6276 (1.94-1.90)
Clashscore	112137	7025 (1.94-1.90)
Ramachandran outliers	110173	6947 (1.94-1.90)
Sidechain outliers	110143	6948 (1.94-1.90)
RSRZ outliers	101464	6332 (1.94-1.90)

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	336	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 12%, green 79%, grey 8%);"></div> <div style="display: flex; justify-content: space-between; font-size: 8px;"> % 79% 12% 8% </div> </div>
1	B	336	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 17%, green 74%, grey 8%);"></div> <div style="display: flex; justify-content: space-between; font-size: 8px;"> % 74% 17% 8% </div> </div>
1	C	336	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, yellow 11%, green 80%, grey 8%);"></div> <div style="display: flex; justify-content: space-between; font-size: 8px;"> 80% 11% 8% </div> </div>
1	D	336	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 17%, green 74%, grey 8%);"></div> <div style="display: flex; justify-content: space-between; font-size: 8px;"> % 74% 17% 8% </div> </div>
1	E	336	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 13%, green 78%, grey 8%);"></div> <div style="display: flex; justify-content: space-between; font-size: 8px;"> % 78% 13% 8% </div> </div>
1	F	336	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, yellow 15%, green 76%, grey 8%);"></div> <div style="display: flex; justify-content: space-between; font-size: 8px;"> 76% 15% 8% </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 15969 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 Iron ABC transporter, periplasmic iron-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	308	Total	C	N	O	S	0	0	0
			2388	1530	419	434	5			
1	B	308	Total	C	N	O	S	0	0	0
			2388	1530	419	434	5			
1	C	308	Total	C	N	O	S	8	1	0
			2396	1535	422	434	5			
1	D	308	Total	C	N	O	S	0	0	0
			2388	1530	419	434	5			
1	E	308	Total	C	N	O	S	0	0	0
			2388	1530	419	434	5			
1	F	308	Total	C	N	O	S	0	0	0
			2388	1530	419	434	5			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	331	HIS	-	EXPRESSION TAG	UNP Q5SHV2
A	332	HIS	-	EXPRESSION TAG	UNP Q5SHV2
A	333	HIS	-	EXPRESSION TAG	UNP Q5SHV2
A	334	HIS	-	EXPRESSION TAG	UNP Q5SHV2
A	335	HIS	-	EXPRESSION TAG	UNP Q5SHV2
A	336	HIS	-	EXPRESSION TAG	UNP Q5SHV2
B	331	HIS	-	EXPRESSION TAG	UNP Q5SHV2
B	332	HIS	-	EXPRESSION TAG	UNP Q5SHV2
B	333	HIS	-	EXPRESSION TAG	UNP Q5SHV2
B	334	HIS	-	EXPRESSION TAG	UNP Q5SHV2
B	335	HIS	-	EXPRESSION TAG	UNP Q5SHV2
B	336	HIS	-	EXPRESSION TAG	UNP Q5SHV2
C	331	HIS	-	EXPRESSION TAG	UNP Q5SHV2
C	332	HIS	-	EXPRESSION TAG	UNP Q5SHV2
C	333	HIS	-	EXPRESSION TAG	UNP Q5SHV2
C	334	HIS	-	EXPRESSION TAG	UNP Q5SHV2
C	335	HIS	-	EXPRESSION TAG	UNP Q5SHV2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	336	HIS	-	EXPRESSION TAG	UNP Q5SHV2
D	331	HIS	-	EXPRESSION TAG	UNP Q5SHV2
D	332	HIS	-	EXPRESSION TAG	UNP Q5SHV2
D	333	HIS	-	EXPRESSION TAG	UNP Q5SHV2
D	334	HIS	-	EXPRESSION TAG	UNP Q5SHV2
D	335	HIS	-	EXPRESSION TAG	UNP Q5SHV2
D	336	HIS	-	EXPRESSION TAG	UNP Q5SHV2
E	331	HIS	-	EXPRESSION TAG	UNP Q5SHV2
E	332	HIS	-	EXPRESSION TAG	UNP Q5SHV2
E	333	HIS	-	EXPRESSION TAG	UNP Q5SHV2
E	334	HIS	-	EXPRESSION TAG	UNP Q5SHV2
E	335	HIS	-	EXPRESSION TAG	UNP Q5SHV2
E	336	HIS	-	EXPRESSION TAG	UNP Q5SHV2
F	331	HIS	-	EXPRESSION TAG	UNP Q5SHV2
F	332	HIS	-	EXPRESSION TAG	UNP Q5SHV2
F	333	HIS	-	EXPRESSION TAG	UNP Q5SHV2
F	334	HIS	-	EXPRESSION TAG	UNP Q5SHV2
F	335	HIS	-	EXPRESSION TAG	UNP Q5SHV2
F	336	HIS	-	EXPRESSION TAG	UNP Q5SHV2

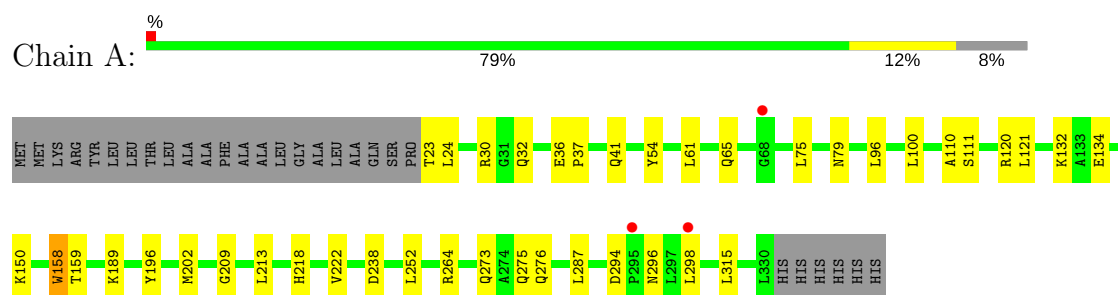
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	293	Total O 293 293	0	0
2	B	248	Total O 248 248	0	0
2	C	331	Total O 331 331	0	0
2	D	252	Total O 252 252	0	0
2	E	252	Total O 252 252	0	0
2	F	257	Total O 257 257	0	0

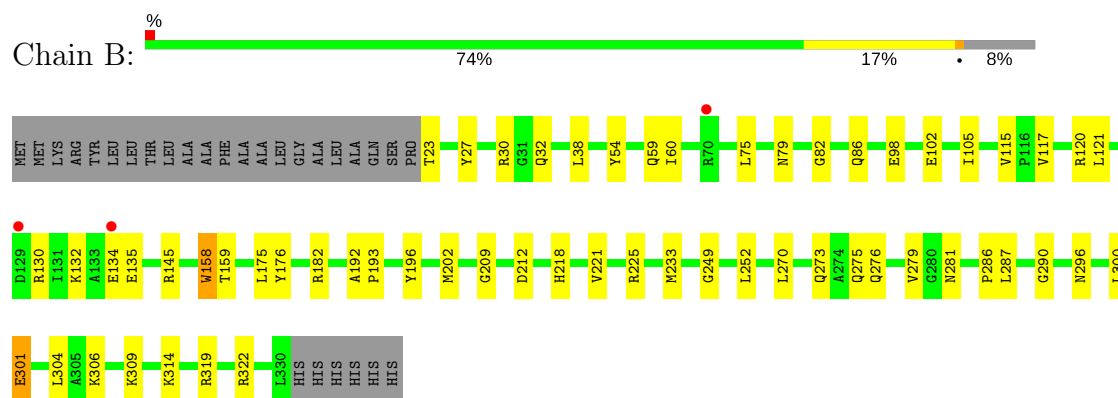
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

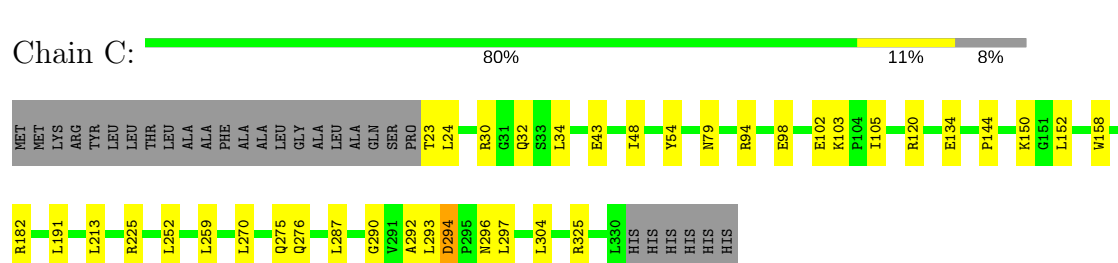
- Molecule 1: Iron ABC transporter, periplasmic iron-binding protein



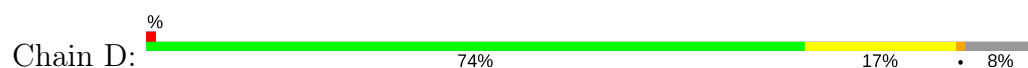
- Molecule 1: Iron ABC transporter, periplasmic iron-binding protein



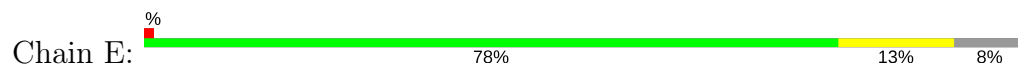
- Molecule 1: Iron ABC transporter, periplasmic iron-binding protein



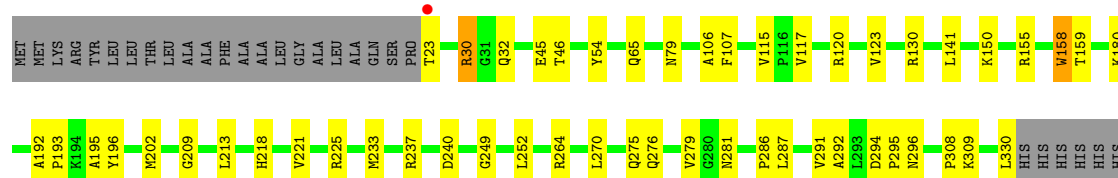
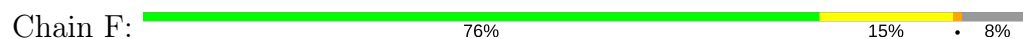
- Molecule 1: Iron ABC transporter, periplasmic iron-binding protein



- Molecule 1: Iron ABC transporter, periplasmic iron-binding protein



- Molecule 1: Iron ABC transporter, periplasmic iron-binding protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	42.12Å 139.35Å 326.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.15 – 1.91 36.15 – 1.91	Depositor EDS
% Data completeness (in resolution range)	94.8 (36.15-1.91) 95.6 (36.15-1.91)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.95 (at 1.91Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_596)	Depositor
R, R_{free}	0.163 , 0.207 0.160 , 0.204	Depositor DCC
R_{free} test set	7181 reflections (5.02%)	DCC
Wilson B-factor (Å ²)	20.8	Xtriage
Anisotropy	0.305	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 56.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	15969	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.71% 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	0.37	0/2438	0.53	0/3311
1	B	0.33	0/2438	0.50	0/3311
1	C	0.39	0/2449	0.55	1/3325 (0.0%)
1	D	0.34	0/2438	0.52	0/3311
1	E	0.32	0/2438	0.48	0/3311
1	F	0.34	0/2438	0.52	2/3311 (0.1%)
All	All	0.35	0/14639	0.52	3/19880 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	30	ARG	NE-CZ-NH2	-5.82	117.39	120.30
1	F	30	ARG	NE-CZ-NH1	5.23	122.91	120.30
1	C	182	ARG	NE-CZ-NH2	-5.12	117.74	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	2388	0	2449	36	0
1	B	2388	0	2449	46	0
1	C	2396	0	2462	27	0
1	D	2388	0	2449	52	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	2388	0	2449	37	0
1	F	2388	0	2449	53	0
2	A	293	0	0	8	0
2	B	248	0	0	8	0
2	C	331	0	0	2	0
2	D	252	0	0	2	0
2	E	252	0	0	6	0
2	F	257	0	0	4	0
All	All	15969	0	14707	244	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 8.

All (244) 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:120:ARG:HH11	1:D:120:ARG:HG3	1.13	1.09
1:F:123:VAL:HG21	1:F:233:MET:HE2	1.43	0.99
1:D:120:ARG:HD2	1:D:219:TYR:CE2	2.01	0.96
1:F:23:THR:HG22	2:F:581:HOH:O	1.67	0.94
1:D:120:ARG:HD3	1:D:284:GLU:OE2	1.70	0.91
1:F:123:VAL:HG21	1:F:233:MET:CE	1.99	0.91
1:C:276:GLN:HE22	1:C:292:ALA:H	1.12	0.90
1:B:98:GLU:O	1:B:102:GLU:HG2	1.71	0.89
1:F:159:THR:HG23	1:F:202:MET:HE1	1.56	0.86
1:F:30:ARG:HD2	1:F:79:ASN:HD22	1.41	0.85
1:D:120:ARG:CG	1:D:120:ARG:HH11	1.91	0.84
1:F:30:ARG:CD	1:F:79:ASN:HD22	1.92	0.82
1:F:221:VAL:HB	1:F:233:MET:HE1	1.60	0.81
1:D:111:SER:OG	1:D:113:THR:HG22	1.83	0.79
1:F:279:VAL:HG22	1:F:286:PRO:HD3	1.65	0.78
1:D:275:GLN:HE22	1:D:287:LEU:H	1.32	0.77
1:D:88:SER:OG	1:D:113:THR:HG21	1.85	0.76
1:B:75:LEU:HD21	1:B:252:LEU:HD23	1.67	0.76
1:B:159:THR:HG23	1:B:202:MET:HE1	1.66	0.76
1:F:276:GLN:HE22	1:F:292:ALA:H	1.31	0.76
1:A:273:GLN:HE21	1:E:276:GLN:HE21	1.34	0.75
1:A:275:GLN:HE22	1:A:287:LEU:H	1.34	0.75
1:A:41:GLN:HE22	1:A:273:GLN:HE22	1.35	0.74
1:E:276:GLN:HE22	1:E:292:ALA:H	1.33	0.74
1:F:275:GLN:HE22	1:F:287:LEU:H	1.35	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:202:MET:SD	2:B:430:HOH:O	2.45	0.74
1:C:275:GLN:HE22	1:C:287:LEU:H	1.33	0.73
1:B:319:ARG:HD2	1:B:322:ARG:NH1	2.05	0.72
1:D:120:ARG:NH1	1:D:120:ARG:HG3	1.95	0.72
1:A:275:GLN:NE2	1:A:287:LEU:H	1.88	0.71
1:D:275:GLN:NE2	1:D:287:LEU:H	1.88	0.71
1:D:120:ARG:HD3	1:D:284:GLU:CD	2.11	0.71
1:D:75:LEU:HD21	1:D:252:LEU:HD23	1.72	0.71
1:F:115:VAL:HG21	1:F:270:LEU:HD11	1.72	0.70
1:E:279:VAL:HG12	1:E:286:PRO:HD3	1.73	0.70
1:F:218:HIS:HA	1:F:233:MET:HE2	1.74	0.69
1:A:273:GLN:HE21	1:E:276:GLN:NE2	1.90	0.69
1:D:75:LEU:CD2	1:D:252:LEU:HD23	2.22	0.69
1:A:75:LEU:HD21	1:A:252:LEU:HD23	1.75	0.68
1:B:30:ARG:HE	1:B:79:ASN:ND2	1.92	0.68
1:D:120:ARG:HD2	1:D:219:TYR:CD2	2.28	0.68
1:C:43:GLU:HG3	1:C:48:ILE:O	1.93	0.67
1:D:41:GLN:HE22	1:D:273:GLN:NE2	1.91	0.67
1:F:30:ARG:HD2	1:F:79:ASN:ND2	2.10	0.66
1:D:115:VAL:HG21	1:D:270:LEU:HD11	1.77	0.66
1:B:275:GLN:HE22	1:B:287:LEU:H	1.44	0.66
1:E:275:GLN:HE22	1:E:287:LEU:H	1.42	0.66
1:A:158:TRP:C	1:A:202:MET:HE1	2.16	0.66
1:B:23:THR:HB	2:B:635:HOH:O	1.96	0.65
1:B:121:LEU:H	1:B:218:HIS:HD2	1.45	0.65
1:A:159:THR:HG23	1:A:202:MET:HE1	1.79	0.65
1:F:130:ARG:HH11	1:F:155:ARG:NH2	1.94	0.65
1:A:134:GLU:H	1:A:134:GLU:CD	1.99	0.64
1:F:275:GLN:NE2	1:F:287:LEU:H	1.96	0.64
1:D:279:VAL:HG22	1:D:286:PRO:HD3	1.80	0.63
1:C:275:GLN:NE2	1:C:287:LEU:H	1.95	0.63
1:A:75:LEU:CD2	1:A:252:LEU:HD23	2.29	0.63
1:A:30:ARG:HE	1:A:79:ASN:ND2	1.97	0.63
1:F:30:ARG:HD3	1:F:79:ASN:HA	1.79	0.63
1:F:130:ARG:NH1	1:F:155:ARG:NH2	2.47	0.62
1:B:75:LEU:CD2	1:B:252:LEU:HD23	2.29	0.62
1:E:302:GLU:O	1:E:306:LYS:HG2	1.98	0.62
1:F:115:VAL:CG2	1:F:270:LEU:HD11	2.30	0.61
1:E:275:GLN:NE2	1:E:287:LEU:H	1.98	0.61
1:A:150:LYS:HG3	1:A:150:LYS:O	1.99	0.61
1:A:132:LYS:N	1:A:132:LYS:HD2	2.16	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:145:ARG:HD3	2:B:572:HOH:O	2.01	0.60
1:B:59:GLN:HB3	2:B:637:HOH:O	2.01	0.60
1:D:158:TRP:HZ3	1:D:195:ALA:N	2.00	0.60
1:A:189:LYS:HE3	2:C:634:HOH:O	2.00	0.60
1:A:264:ARG:NH2	2:A:630:HOH:O	2.35	0.60
1:C:294:ASP:H	1:C:297:LEU:HD12	1.65	0.59
1:C:276:GLN:NE2	1:C:292:ALA:H	1.94	0.59
1:D:158:TRP:CZ3	1:D:195:ALA:HB2	2.37	0.59
1:A:276:GLN:HE21	1:E:281:ASN:HD21	1.49	0.59
1:F:240:ASP:O	1:F:308:PRO:HG3	2.03	0.59
1:F:123:VAL:HG21	1:F:233:MET:HE3	1.85	0.59
1:B:196:TYR:HD2	1:B:202:MET:HE2	1.67	0.58
1:C:276:GLN:HE22	1:C:292:ALA:N	1.94	0.58
1:C:252:LEU:HD21	1:C:259:LEU:CD1	2.33	0.58
1:D:120:ARG:CD	1:D:219:TYR:CE2	2.84	0.57
1:B:275:GLN:NE2	1:B:287:LEU:H	2.03	0.57
1:A:30:ARG:HE	1:A:79:ASN:HD22	1.50	0.57
1:E:307:SER:HB2	2:E:435:HOH:O	2.05	0.57
1:F:150:LYS:HG3	1:F:150:LYS:O	2.04	0.57
1:F:158:TRP:CZ3	1:F:195:ALA:HB2	2.40	0.56
1:B:319:ARG:HD2	1:B:322:ARG:HH11	1.70	0.56
1:B:301:GLU:H	1:B:301:GLU:CD	2.09	0.56
1:E:121:LEU:H	1:E:218:HIS:HD2	1.54	0.56
1:C:252:LEU:HD21	1:C:259:LEU:HD11	1.88	0.56
1:C:30:ARG:HE	1:C:79:ASN:ND2	2.04	0.55
1:D:132:LYS:HD2	1:D:132:LYS:N	2.21	0.55
1:B:158:TRP:C	1:B:202:MET:HE1	2.26	0.55
1:B:30:ARG:HE	1:B:79:ASN:HD22	1.52	0.55
1:B:132:LYS:O	1:B:135:GLU:HG2	2.07	0.55
1:B:132:LYS:HG3	1:B:134:GLU:OE1	2.07	0.55
1:F:209:GLY:HA2	2:F:445:HOH:O	2.05	0.55
1:F:30:ARG:HD3	1:F:79:ASN:HD22	1.70	0.55
1:F:192:ALA:N	1:F:193:PRO:HD3	2.22	0.55
1:A:159:THR:HG23	1:A:202:MET:CE	2.37	0.54
1:F:159:THR:HG23	1:F:202:MET:CE	2.33	0.54
1:A:196:TYR:HD2	1:A:202:MET:HE2	1.72	0.54
1:B:309:LYS:HE3	2:B:448:HOH:O	2.08	0.54
1:C:293:LEU:O	1:D:37:PRO:HB3	2.08	0.54
1:E:61:LEU:O	1:E:65:GLN:HG2	2.08	0.53
1:C:290:GLY:HA2	1:D:45:GLU:OE2	2.09	0.53
1:C:134:GLU:CD	1:C:134:GLU:H	2.11	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:120:ARG:NH1	1:D:120:ARG:CG	2.60	0.53
1:E:179:GLU:HG2	2:E:573:HOH:O	2.09	0.53
1:C:30:ARG:HE	1:C:79:ASN:HD22	1.58	0.52
1:F:218:HIS:HA	1:F:233:MET:CE	2.38	0.52
1:D:30:ARG:HE	1:D:79:ASN:ND2	2.07	0.52
1:F:221:VAL:HB	1:F:233:MET:CE	2.37	0.52
1:B:115:VAL:HG21	1:B:270:LEU:HD11	1.91	0.52
1:B:159:THR:HG23	1:B:202:MET:CE	2.38	0.52
1:D:158:TRP:CZ2	1:D:193:PRO:HG2	2.45	0.52
1:C:150:LYS:HG3	1:C:150:LYS:O	2.10	0.51
1:F:158:TRP:HZ3	1:F:195:ALA:HB2	1.76	0.51
1:D:225:ARG:HH11	1:D:296:ASN:ND2	2.09	0.51
1:F:308:PRO:HD2	2:F:469:HOH:O	2.11	0.50
1:E:279:VAL:HG11	1:E:286:PRO:HG3	1.94	0.50
1:F:30:ARG:CD	1:F:79:ASN:ND2	2.67	0.50
1:D:213:LEU:C	1:D:213:LEU:HD12	2.31	0.50
1:E:34:LEU:HD23	1:E:34:LEU:C	2.32	0.50
1:A:121:LEU:H	1:A:218:HIS:HD2	1.59	0.50
1:A:23:THR:N	2:A:495:HOH:O	2.45	0.50
1:F:309:LYS:N	1:F:309:LYS:HD2	2.27	0.49
1:D:115:VAL:CG2	1:D:270:LEU:HD11	2.41	0.49
1:C:103:LYS:HB2	1:C:270:LEU:HD23	1.93	0.49
1:F:225:ARG:HH11	1:F:296:ASN:ND2	2.11	0.49
1:C:94:ARG:HD2	1:C:252:LEU:HG	1.92	0.49
1:D:41:GLN:HE22	1:D:273:GLN:HE22	1.60	0.49
1:F:158:TRP:C	1:F:202:MET:HE1	2.34	0.48
1:F:221:VAL:HG21	1:F:233:MET:HE3	1.95	0.48
1:F:130:ARG:NH1	1:F:155:ARG:HH22	2.11	0.48
1:F:202:MET:HB2	1:F:202:MET:HE3	1.44	0.48
1:A:213:LEU:C	1:A:213:LEU:HD12	2.34	0.48
1:E:279:VAL:CG1	1:E:286:PRO:HD3	2.42	0.48
1:A:23:THR:HG22	1:A:24:LEU:N	2.28	0.48
1:A:61:LEU:O	1:A:65:GLN:HG2	2.14	0.48
1:D:36:GLU:HG2	1:D:40:LYS:HE2	1.96	0.48
1:F:46:THR:HG22	1:F:264:ARG:HH21	1.78	0.48
1:D:103:LYS:NZ	1:D:267:THR:HG23	2.30	0.47
1:D:63:ALA:O	1:D:66:GLU:HG2	2.14	0.47
1:E:23:THR:N	2:E:558:HOH:O	2.46	0.47
1:B:121:LEU:H	1:B:218:HIS:CD2	2.29	0.47
1:B:182:ARG:NH2	2:B:620:HOH:O	2.48	0.47
1:E:121:LEU:H	1:E:218:HIS:CD2	2.31	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:209:GLY:HA2	2:E:452:HOH:O	2.14	0.47
1:B:279:VAL:HG22	1:B:286:PRO:HD3	1.97	0.47
1:D:75:LEU:HD23	1:D:252:LEU:HD23	1.97	0.47
1:B:290:GLY:HA2	1:F:45:GLU:OE2	2.15	0.47
1:E:238:ASP:OD1	1:E:306:LYS:HD3	2.15	0.46
1:D:158:TRP:CH2	1:D:193:PRO:HG2	2.50	0.46
1:E:39:VAL:O	1:E:43:GLU:HG3	2.16	0.46
1:C:144:PRO:HA	1:C:191:LEU:HD11	1.96	0.46
1:A:222:VAL:HG22	1:A:296:ASN:HB3	1.98	0.46
1:E:225:ARG:HH11	1:E:296:ASN:ND2	2.14	0.46
1:B:202:MET:HE3	1:B:202:MET:HB2	1.36	0.46
1:B:27:TYR:HB3	1:B:60:ILE:HD12	1.97	0.46
1:A:111:SER:N	2:A:539:HOH:O	2.48	0.46
1:D:138:GLU:O	1:D:236:PHE:HB3	2.15	0.46
1:E:225:ARG:HH11	1:E:296:ASN:HD22	1.63	0.45
1:A:121:LEU:H	1:A:218:HIS:CD2	2.34	0.45
1:A:32:GLN:HA	1:A:54:TYR:CE1	2.51	0.45
1:C:325:ARG:HD3	2:C:523:HOH:O	2.16	0.45
1:E:74:ASP:HA	1:E:256:SER:HB2	1.97	0.45
1:D:292:ALA:O	2:D:433:HOH:O	2.21	0.45
1:A:96:LEU:HD13	1:A:100:LEU:HD23	1.98	0.45
1:D:132:LYS:HB3	1:D:134:GLU:OE2	2.17	0.45
1:D:43:GLU:HG2	1:D:48:ILE:O	2.16	0.45
1:A:110:ALA:HB3	2:A:539:HOH:O	2.16	0.45
1:D:158:TRP:CZ3	1:D:195:ALA:CA	3.00	0.45
1:E:318:ASP:HB2	2:E:423:HOH:O	2.17	0.45
1:F:330:LEU:HA	1:F:330:LEU:HD12	1.80	0.45
1:B:225:ARG:HH11	1:B:296:ASN:HD22	1.65	0.45
1:D:32:GLN:HA	1:D:54:TYR:CE1	2.52	0.45
1:E:23:THR:HG22	2:E:432:HOH:O	2.17	0.45
1:B:281:ASN:HD22	1:B:281:ASN:H	1.64	0.45
1:F:286:PRO:HG3	1:F:291:VAL:HG21	1.99	0.45
1:A:264:ARG:HD2	2:A:485:HOH:O	2.18	0.44
1:B:130:ARG:HD2	1:B:212:ASP:OD1	2.17	0.44
1:B:32:GLN:HA	1:B:54:TYR:CE1	2.52	0.44
1:B:105:ILE:HD13	1:B:287:LEU:O	2.18	0.44
1:C:34:LEU:C	1:C:34:LEU:HD23	2.38	0.44
1:F:141:LEU:HD13	1:F:180:LYS:HE3	2.00	0.44
1:D:27:TYR:HB3	1:D:60:ILE:HD12	1.99	0.44
1:D:119:VAL:HG13	1:D:285:TYR:HB2	2.00	0.44
1:F:32:GLN:HA	1:F:54:TYR:CE1	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:172:MET:HE2	1:D:181:THR:HG23	2.00	0.43
1:C:98:GLU:O	1:C:102:GLU:HG3	2.17	0.43
1:B:132:LYS:HD2	1:B:132:LYS:HA	1.72	0.43
1:B:117:VAL:HG23	1:B:249:GLY:HA2	2.00	0.43
1:F:213:LEU:C	1:F:213:LEU:HD12	2.38	0.43
1:B:209:GLY:HA2	2:B:432:HOH:O	2.17	0.43
1:E:175:LEU:HD11	1:E:308:PRO:HB3	2.00	0.43
1:B:175:LEU:HD23	1:B:176:TYR:CZ	2.54	0.43
1:C:23:THR:HG22	1:C:24:LEU:N	2.34	0.43
1:D:23:THR:N	2:D:480:HOH:O	2.51	0.43
1:A:36:GLU:HB3	1:A:37:PRO:HD3	2.01	0.43
1:D:127:ASN:HA	1:D:128:PRO:HD3	1.88	0.43
1:E:99:THR:HG23	1:E:100:LEU:N	2.33	0.43
1:C:32:GLN:HA	1:C:54:TYR:CE1	2.54	0.43
1:E:324:LEU:HB3	1:E:330:LEU:HB2	2.01	0.43
1:B:38:LEU:HD23	1:B:38:LEU:HA	1.88	0.43
1:C:150:LYS:HG2	1:C:152:LEU:HD11	2.00	0.43
1:F:237:ARG:HD2	2:F:540:HOH:O	2.19	0.43
1:C:213:LEU:HD12	1:C:213:LEU:C	2.38	0.42
1:A:315:LEU:N	2:A:554:HOH:O	2.51	0.42
1:C:225:ARG:HD3	1:C:296:ASN:ND2	2.34	0.42
1:F:196:TYR:HD2	1:F:202:MET:HE2	1.84	0.42
1:E:105:ILE:HG12	1:E:287:LEU:O	2.19	0.42
1:F:106:ALA:O	1:F:107:PHE:HB2	2.19	0.42
1:A:209:GLY:HA2	2:A:464:HOH:O	2.19	0.42
1:F:117:VAL:HG23	1:F:249:GLY:HA2	2.02	0.42
1:E:29:GLY:HA2	1:E:55:SER:O	2.19	0.42
1:F:46:THR:HG22	1:F:264:ARG:NH2	2.35	0.42
1:E:110:ALA:O	1:E:112:ARG:HG2	2.19	0.42
1:F:294:ASP:HA	1:F:295:PRO:HD3	1.96	0.42
1:E:276:GLN:HA	1:E:279:VAL:HG22	2.02	0.42
1:E:279:VAL:HG23	1:E:280:GLY:N	2.35	0.42
1:A:202:MET:HE3	1:A:202:MET:HB2	1.60	0.41
1:E:32:GLN:HA	1:E:54:TYR:CE1	2.54	0.41
1:D:144:PRO:HA	1:D:191:LEU:HD11	2.02	0.41
1:F:159:THR:CG2	1:F:202:MET:HE1	2.40	0.41
1:A:315:LEU:HB2	2:A:554:HOH:O	2.21	0.41
1:D:233:MET:HG2	1:D:233:MET:O	2.20	0.41
1:E:175:LEU:HD12	1:E:241:ALA:HB2	2.03	0.41
1:E:71:SER:HA	1:E:72:PRO:HD3	1.81	0.41
1:B:276:GLN:HE21	1:F:281:ASN:HD21	1.69	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:299:PRO:HG2	1:D:302:GLU:OE2	2.21	0.41
1:D:30:ARG:HE	1:D:79:ASN:HD22	1.67	0.41
1:B:300:LEU:O	1:B:304:LEU:HG	2.20	0.41
1:B:306:LYS:HD3	1:B:306:LYS:HA	1.88	0.41
1:C:105:ILE:CD1	1:C:304:LEU:HD11	2.51	0.41
1:D:281:ASN:HD22	1:D:281:ASN:H	1.69	0.41
1:B:221:VAL:HG11	1:B:233:MET:HB2	2.03	0.40
1:B:82:GLY:O	1:B:86:GLN:HG3	2.21	0.40
1:F:115:VAL:HG21	1:F:270:LEU:CD1	2.45	0.40
1:B:192:ALA:N	1:B:193:PRO:HD3	2.36	0.40
1:D:103:LYS:N	1:D:104:PRO:CD	2.84	0.40
1:E:301:GLU:H	1:E:301:GLU:HG3	1.51	0.40
1:B:314:LYS:HE3	2:B:592:HOH:O	2.21	0.40
1:D:132:LYS:O	1:D:135:GLU:HB2	2.21	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	306/336 (91%)	299 (98%)	7 (2%)	0	100	100
1	B	306/336 (91%)	304 (99%)	2 (1%)	0	100	100
1	C	307/336 (91%)	305 (99%)	2 (1%)	0	100	100
1	D	306/336 (91%)	304 (99%)	2 (1%)	0	100	100
1	E	306/336 (91%)	305 (100%)	1 (0%)	0	100	100
1	F	306/336 (91%)	304 (99%)	2 (1%)	0	100	100
All	All	1837/2016 (91%)	1821 (99%)	16 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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	247/268 (92%)	242 (98%)	5 (2%)	60	54
1	B	247/268 (92%)	243 (98%)	4 (2%)	68	63
1	C	248/268 (92%)	245 (99%)	3 (1%)	75	74
1	D	247/268 (92%)	242 (98%)	5 (2%)	60	54
1	E	247/268 (92%)	244 (99%)	3 (1%)	75	74
1	F	247/268 (92%)	243 (98%)	4 (2%)	68	63
All	All	1483/1608 (92%)	1459 (98%)	24 (2%)	68	63

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	120	ARG
1	A	158	TRP
1	A	238	ASP
1	A	294	ASP
1	A	298	LEU
1	B	120	ARG
1	B	158	TRP
1	B	273	GLN
1	B	301	GLU
1	C	120	ARG
1	C	158	TRP
1	C	294	ASP
1	D	120	ARG
1	D	158	TRP
1	D	273	GLN
1	D	300	LEU
1	D	323	LEU
1	E	120	ARG
1	E	158	TRP
1	E	301	GLU
1	F	65	GLN
1	F	120	ARG

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Mol	Chain	Res	Type
1	F	158	TRP
1	F	252	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	GLN
1	A	65	GLN
1	A	79	ASN
1	A	218	HIS
1	A	275	GLN
1	A	281	ASN
1	B	65	GLN
1	B	79	ASN
1	B	218	HIS
1	B	273	GLN
1	B	275	GLN
1	B	281	ASN
1	B	296	ASN
1	C	79	ASN
1	C	275	GLN
1	C	276	GLN
1	C	296	ASN
1	D	79	ASN
1	D	273	GLN
1	D	275	GLN
1	D	281	ASN
1	D	296	ASN
1	E	32	GLN
1	E	65	GLN
1	E	199	ASN
1	E	218	HIS
1	E	275	GLN
1	E	276	GLN
1	E	281	ASN
1	E	296	ASN
1	F	65	GLN
1	F	79	ASN
1	F	275	GLN
1	F	276	GLN
1	F	281	ASN
1	F	296	ASN

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	308/336 (91%)	-0.19	3 (0%) 82 84	9, 17, 42, 59	0
1	B	308/336 (91%)	-0.17	3 (0%) 82 84	11, 23, 46, 66	0
1	C	308/336 (91%)	-0.36	0 100 100	7, 15, 33, 49	0
1	D	308/336 (91%)	-0.12	5 (1%) 72 76	12, 23, 48, 64	0
1	E	308/336 (91%)	-0.18	3 (0%) 82 84	12, 23, 47, 65	0
1	F	308/336 (91%)	-0.22	1 (0%) 93 94	12, 22, 45, 58	0
All	All	1848/2016 (91%)	-0.21	15 (0%) 86 88	7, 21, 45, 66	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	326	GLU	3.5
1	A	298	LEU	3.0
1	F	23	THR	2.8
1	E	230	ARG	2.7
1	D	99	THR	2.5
1	A	68	GLY	2.4
1	B	70	ARG	2.3
1	D	297	LEU	2.3
1	D	98	GLU	2.2
1	A	295	PRO	2.2
1	B	134	GLU	2.1
1	D	23	THR	2.1
1	D	158	TRP	2.1
1	E	99	THR	2.0
1	B	129	ASP	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.