



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2023 – 11:52 AM EDT

PDB ID : 2RG7
Title : Apo- Crystal Structure of a Periplasmic Heme Binding Protein from Shigella dysenteriae
Authors : Ho, W.W.; Li, H.
Deposited on : 2007-10-02
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 : 2.35
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) : 2.35

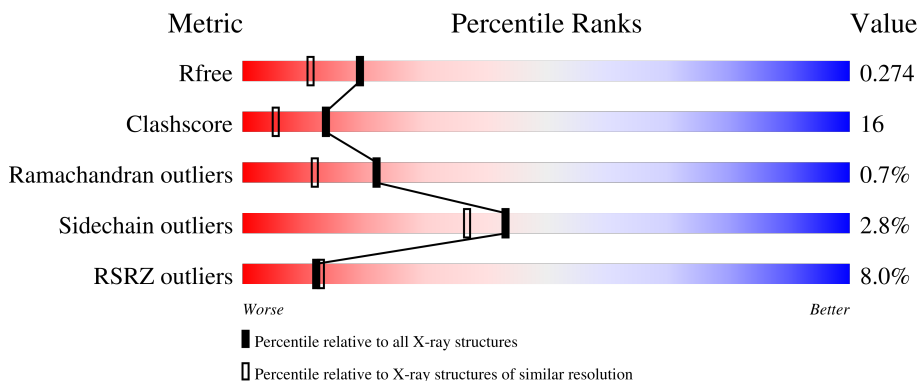
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


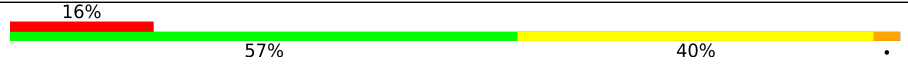
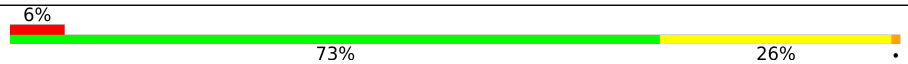

The reported resolution of this entry is 2.05 Å.

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	256	
1	B	256	
1	C	256	
1	D	256	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7982 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 Periplasmic HEME binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	255	1934	1216	344	369	5	0	0	0
1	B	255	1934	1216	344	369	5	0	0	0
1	C	255	1934	1216	344	369	5	0	0	0
1	D	255	1934	1216	344	369	5	0	0	0

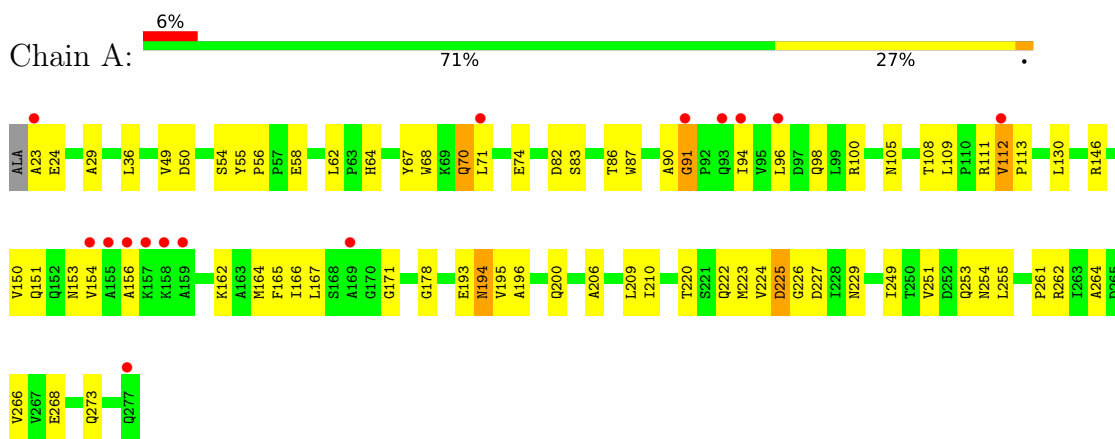
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	73	Total	O	0	0
			73	73		
2	B	54	Total	O	0	0
			54	54		
2	C	58	Total	O	0	0
			58	58		
2	D	61	Total	O	0	0
			61	61		

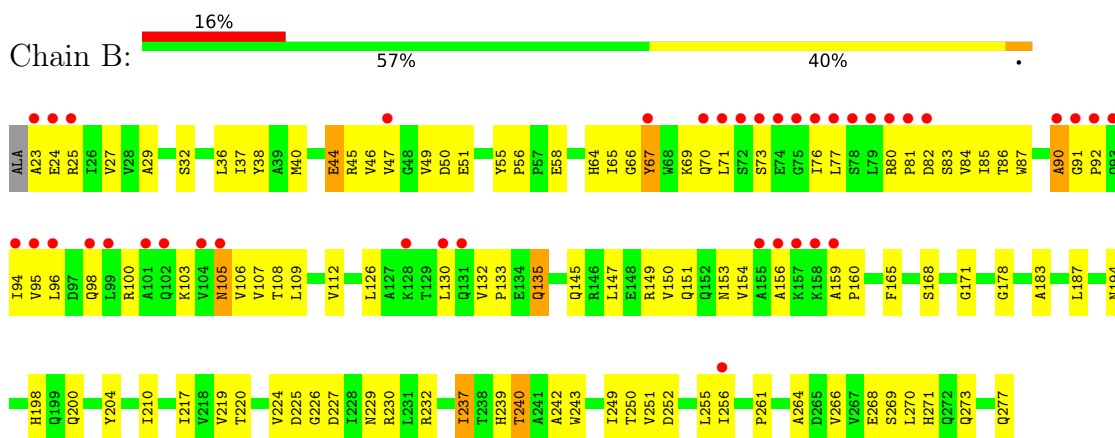
3 Residue-property plots [i](#)

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.

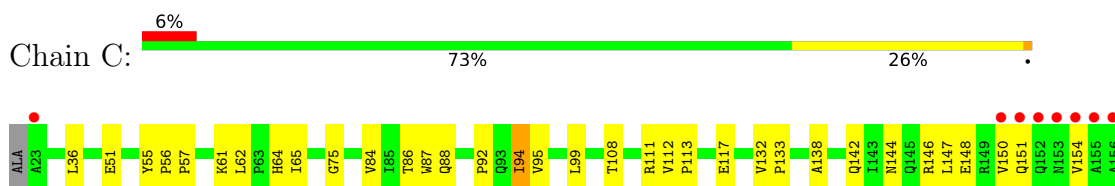
- Molecule 1: Periplasmic HEME binding protein

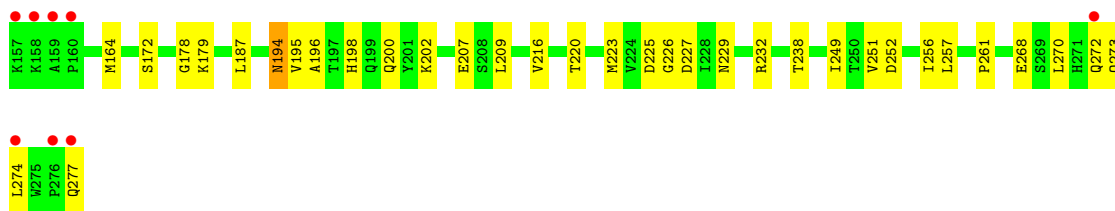


- Molecule 1: Periplasmic HEME binding protein

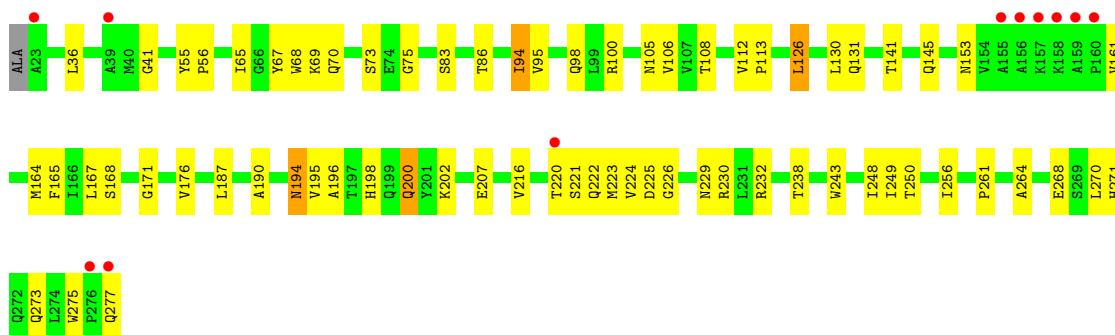
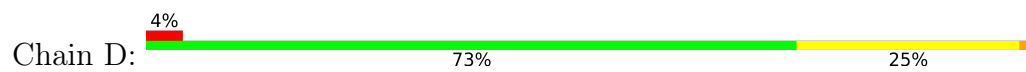


- Molecule 1: Periplasmic HEME binding protein





- Molecule 1: Periplasmic HEME binding protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	54.82Å 73.28Å 72.38Å 71.44° 77.80° 89.90°	Depositor
Resolution (Å)	35.90 – 2.05 35.90 – 2.05	Depositor EDS
% Data completeness (in resolution range)	97.4 (35.90-2.05) 97.5 (35.90-2.05)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.58 (at 2.05Å)	Xtrriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.232 , 0.274 0.232 , 0.274	Depositor DCC
R_{free} test set	6402 reflections (10.10%)	wwPDB-VP
Wilson B-factor (Å ²)	38.3	Xtrriage
Anisotropy	0.463	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 50.7	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	0.95	EDS
Total number of atoms	7982	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.62% 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	0.33	0/1965	0.59	0/2676
1	B	0.30	0/1965	0.59	0/2676
1	C	0.33	0/1965	0.57	0/2676
1	D	0.33	0/1965	0.59	0/2676
All	All	0.32	0/7860	0.58	0/10704

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	1934	0	1978	52	0
1	B	1934	0	1978	98	0
1	C	1934	0	1978	54	0
1	D	1934	0	1978	54	0
2	A	73	0	0	4	0
2	B	54	0	0	2	0
2	C	58	0	0	2	0
2	D	61	0	0	3	0
All	All	7982	0	7912	249	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 16.

All (249) 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:A:112:VAL:HG13	1:A:113:PRO:HD3	1.31	1.06
1:B:94:ILE:HD12	1:B:94:ILE:H	1.31	0.96
1:D:67:TYR:HB2	1:D:70:GLN:HG2	1.59	0.84
1:A:220:THR:H	1:A:223:MET:HE3	1.43	0.84
1:A:67:TYR:HB3	1:A:70:GLN:HG3	1.58	0.83
1:C:207:GLU:HG2	1:D:98:GLN:OE1	1.77	0.83
1:B:240:THR:HG22	1:B:243:TRP:H	1.44	0.83
1:A:255:LEU:HD12	1:A:266:VAL:HG22	1.61	0.83
1:B:23:ALA:N	1:B:83:SER:HG	1.76	0.82
1:D:229:ASN:ND2	1:D:232:ARG:HE	1.78	0.81
1:C:138:ALA:O	1:C:142:GLN:HG3	1.81	0.80
1:D:220:THR:HG23	1:D:223:MET:H	1.46	0.78
1:B:36:LEU:HD23	1:B:261:PRO:HD3	1.67	0.76
1:C:94:ILE:HD13	1:C:94:ILE:O	1.88	0.74
1:C:87:TRP:CE3	1:C:111:ARG:HB3	2.23	0.73
1:C:92:PRO:HG2	1:C:95:VAL:HG23	1.68	0.73
1:A:94:ILE:O	1:A:98:GLN:HG3	1.89	0.72
1:A:150:VAL:O	1:A:154:VAL:HG23	1.89	0.72
1:D:94:ILE:HD13	1:D:94:ILE:O	1.88	0.72
1:D:243:TRP:HA	1:D:248:ILE:HD11	1.72	0.72
1:B:150:VAL:O	1:B:154:VAL:HG23	1.90	0.71
1:B:24:GLU:OE1	1:B:25:ARG:HG3	1.90	0.71
1:C:187:LEU:HD23	1:C:270:LEU:HD21	1.71	0.71
1:B:109:LEU:HD21	1:B:126:LEU:HG	1.73	0.71
1:C:249:ILE:HG23	1:C:273:GLN:OE1	1.91	0.71
1:C:200:GLN:HB3	2:C:333:HOH:O	1.90	0.70
1:B:67:TYR:CE2	1:B:69:LYS:HB3	2.26	0.69
1:A:68:TRP:HA	1:A:71:LEU:CD1	2.23	0.68
1:B:252:ASP:HB3	1:B:255:LEU:HD23	1.76	0.68
1:A:227:ASP:OD2	1:A:229:ASN:HB2	1.94	0.68
1:A:249:ILE:HG23	1:A:273:GLN:OE1	1.92	0.68
1:C:207:GLU:HG2	1:D:98:GLN:CD	2.14	0.68
1:B:58:GLU:CD	1:B:58:GLU:H	1.98	0.67
1:A:220:THR:HG21	1:A:253:GLN:HG3	1.75	0.67
1:B:237:ILE:O	1:B:240:THR:HB	1.94	0.67
1:A:67:TYR:HB3	1:A:70:GLN:CG	2.25	0.67
1:B:73:SER:O	1:B:77:LEU:HG	1.95	0.67
1:D:264:ALA:O	1:D:268:GLU:HG3	1.95	0.67
1:C:179:LYS:HE2	1:C:196:ALA:O	1.95	0.66
1:B:94:ILE:HG22	1:B:98:GLN:HE21	1.61	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:225:ASP:OD1	1:C:226:GLY:N	2.29	0.65
1:B:71:LEU:HD11	1:B:76:ILE:HD11	1.78	0.65
1:B:255:LEU:HB3	1:B:266:VAL:HG21	1.78	0.65
1:B:103:LYS:HE2	1:B:103:LYS:HA	1.79	0.64
1:A:162:LYS:HE3	1:A:193:GLU:OE1	1.97	0.64
1:B:149:ARG:HD2	2:C:280:HOH:O	1.96	0.63
1:D:194:ASN:ND2	1:D:196:ALA:H	1.95	0.63
1:D:187:LEU:HD23	1:D:270:LEU:HD21	1.79	0.62
1:A:255:LEU:HD12	1:A:266:VAL:CG2	2.28	0.62
1:D:36:LEU:HD23	1:D:261:PRO:HG3	1.82	0.62
1:B:92:PRO:HB2	1:B:95:VAL:HG23	1.81	0.62
1:D:94:ILE:O	1:D:98:GLN:HG3	2.00	0.62
1:D:225:ASP:CG	1:D:226:GLY:H	2.03	0.62
1:A:68:TRP:HA	1:A:71:LEU:HD11	1.83	0.60
1:B:132:VAL:HB	1:B:135:GLN:NE2	2.15	0.60
1:B:165:PHE:CZ	1:B:256:ILE:HD13	2.36	0.60
1:B:271:HIS:HE1	1:C:117:GLU:OE2	1.85	0.60
2:A:325:HOH:O	1:C:172:SER:HB3	2.02	0.59
1:B:256:ILE:HG12	1:B:256:ILE:O	2.02	0.59
1:B:187:LEU:HD23	1:B:270:LEU:HD21	1.84	0.59
1:B:264:ALA:O	1:B:268:GLU:HG3	2.03	0.59
1:C:194:ASN:ND2	1:C:196:ALA:H	2.01	0.58
1:B:168:SER:O	1:D:171:GLY:HA2	2.02	0.58
1:A:58:GLU:CD	1:A:58:GLU:H	2.07	0.58
1:A:222:GLN:OE1	1:A:253:GLN:HB2	2.03	0.58
1:B:44:GLU:OE1	1:B:45:ARG:HG3	2.02	0.58
1:B:94:ILE:O	1:B:98:GLN:HG3	2.03	0.58
1:D:194:ASN:HD22	1:D:195:VAL:N	2.02	0.58
1:C:198:HIS:HB2	1:C:202:LYS:HE3	1.85	0.57
1:B:94:ILE:H	1:B:94:ILE:CD1	2.08	0.57
1:C:65:ILE:HG21	1:C:75:GLY:HA3	1.86	0.57
1:B:91:GLY:HA2	1:B:92:PRO:C	2.25	0.57
1:B:70:GLN:HA	1:B:92:PRO:HD3	1.86	0.57
1:A:68:TRP:O	1:A:91:GLY:N	2.28	0.57
1:B:255:LEU:HB3	1:B:266:VAL:CG2	2.33	0.56
1:B:249:ILE:HG23	1:B:273:GLN:OE1	2.04	0.56
1:A:62:LEU:O	1:A:64:HIS:HD2	1.89	0.56
1:A:90:ALA:O	1:A:91:GLY:O	2.24	0.56
1:A:220:THR:H	1:A:223:MET:CE	2.17	0.56
1:C:55:TYR:CG	1:C:56:PRO:HA	2.40	0.56
1:D:112:VAL:HG13	1:D:113:PRO:HA	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:ILE:O	1:A:223:MET:HE1	2.06	0.56
1:B:237:ILE:O	1:B:237:ILE:HD13	2.05	0.56
1:C:268:GLU:O	1:C:272:GLN:HG3	2.06	0.56
1:D:220:THR:HG23	1:D:223:MET:N	2.19	0.56
1:B:37:ILE:HG12	1:B:130:LEU:HD11	1.89	0.55
1:B:49:VAL:O	1:B:64:HIS:HA	2.06	0.55
1:D:243:TRP:HA	1:D:248:ILE:CD1	2.35	0.55
1:C:36:LEU:HD23	1:C:261:PRO:HG3	1.89	0.54
1:B:108:THR:C	1:B:109:LEU:HD12	2.28	0.54
1:B:24:GLU:CD	1:B:25:ARG:HG3	2.27	0.54
1:B:67:TYR:HD2	1:B:70:GLN:HG3	1.73	0.54
1:D:95:VAL:HA	1:D:98:GLN:HE21	1.73	0.54
2:A:318:HOH:O	1:C:232:ARG:HD3	2.07	0.53
1:B:94:ILE:HG22	1:B:98:GLN:NE2	2.23	0.53
1:D:131:GLN:OE1	1:D:131:GLN:HA	2.08	0.53
1:B:38:TYR:CE1	1:B:46:VAL:HG21	2.44	0.53
1:A:164:MET:SD	1:A:209:LEU:HD11	2.49	0.53
1:B:132:VAL:N	1:B:133:PRO:HD3	2.23	0.53
1:C:194:ASN:HD22	1:C:195:VAL:N	2.07	0.53
1:D:141:THR:O	1:D:145:GLN:HG3	2.09	0.53
1:D:65:ILE:HG21	1:D:75:GLY:HA3	1.91	0.52
1:C:150:VAL:O	1:C:154:VAL:HG23	2.09	0.52
1:C:207:GLU:CD	1:C:207:GLU:H	2.13	0.52
1:B:227:ASP:OD2	1:B:229:ASN:HB2	2.09	0.52
1:A:220:THR:HA	1:A:251:VAL:O	2.10	0.52
1:B:219:VAL:O	1:B:250:THR:HA	2.10	0.52
1:A:87:TRP:CE3	1:A:111:ARG:HB2	2.45	0.51
1:B:220:THR:HA	1:B:251:VAL:O	2.09	0.51
1:B:29:ALA:HA	1:B:50:ASP:HB3	1.93	0.51
1:B:153:ASN:O	1:B:156:ALA:HB3	2.11	0.51
1:D:225:ASP:CG	1:D:226:GLY:N	2.64	0.51
1:B:67:TYR:CD2	1:B:69:LYS:HB3	2.45	0.51
1:C:178:GLY:HA3	1:C:200:GLN:O	2.11	0.51
1:A:194:ASN:HD22	1:A:195:VAL:N	2.09	0.50
1:A:86:THR:O	1:A:108:THR:HA	2.12	0.50
1:C:198:HIS:HB2	1:C:202:LYS:CE	2.41	0.50
1:A:55:TYR:CG	1:A:56:PRO:HA	2.46	0.50
1:A:206:ALA:O	1:A:210:ILE:HG13	2.11	0.50
1:C:87:TRP:CD2	1:C:111:ARG:HB3	2.46	0.50
1:C:256:ILE:HD12	1:C:257:LEU:HG	1.94	0.50
1:C:147:LEU:O	1:C:151:GLN:HG3	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:100:ARG:NH1	1:D:106:VAL:HB	2.26	0.50
1:D:249:ILE:HG23	1:D:273:GLN:OE1	2.11	0.50
1:A:54:SER:OG	1:A:254:ASN:ND2	2.45	0.50
1:D:216:VAL:HG11	1:D:249:ILE:CD1	2.42	0.50
1:D:94:ILE:HD13	1:D:98:GLN:HG3	1.95	0.49
1:A:36:LEU:HD23	1:A:261:PRO:HD3	1.95	0.49
1:B:90:ALA:HB3	1:B:96:LEU:HD11	1.93	0.49
1:C:92:PRO:HG2	1:C:95:VAL:CG2	2.41	0.49
1:D:83:SER:HB3	1:D:105:ASN:HB3	1.94	0.49
1:B:38:TYR:HE1	1:B:46:VAL:HG21	1.78	0.49
1:B:73:SER:HB3	1:B:95:VAL:HG13	1.95	0.49
1:B:165:PHE:CE1	1:B:256:ILE:HD13	2.48	0.49
1:B:67:TYR:O	1:B:71:LEU:N	2.45	0.48
1:B:90:ALA:HB3	1:B:96:LEU:CD1	2.43	0.48
1:D:225:ASP:HB3	1:D:230:ARG:NH1	2.27	0.48
1:B:103:LYS:HE2	1:B:103:LYS:CA	2.41	0.48
1:A:264:ALA:O	1:A:268:GLU:HG3	2.14	0.48
1:D:68:TRP:CZ2	1:D:69:LYS:HE2	2.48	0.48
1:D:198:HIS:CG	1:D:202:LYS:HG3	2.48	0.48
1:B:94:ILE:HD12	1:B:94:ILE:N	2.14	0.48
2:A:344:HOH:O	1:C:51:GLU:HG2	2.13	0.48
1:B:277:GLN:NE2	1:C:148:GLU:HG3	2.29	0.48
1:A:194:ASN:ND2	1:A:196:ALA:H	2.11	0.47
1:D:224:VAL:O	1:D:225:ASP:HB3	2.15	0.47
1:C:112:VAL:HG13	1:C:113:PRO:HA	1.97	0.47
1:B:85:ILE:HA	1:B:107:VAL:O	2.15	0.47
1:B:29:ALA:HB3	1:B:86:THR:HB	1.96	0.47
1:C:216:VAL:HG11	1:C:249:ILE:CD1	2.45	0.47
1:D:190:ALA:O	1:D:271:HIS:HD2	1.98	0.46
1:D:256:ILE:C	1:D:256:ILE:HD12	2.34	0.46
1:B:210:ILE:HD11	1:B:239:HIS:HB2	1.98	0.46
1:A:96:LEU:O	1:A:100:ARG:HG3	2.15	0.46
1:C:132:VAL:N	1:C:133:PRO:CD	2.79	0.46
1:C:146:ARG:HH21	1:C:268:GLU:CD	2.18	0.46
1:C:220:THR:HA	1:C:251:VAL:O	2.14	0.46
1:D:200:GLN:HB3	2:D:322:HOH:O	2.14	0.46
1:C:84:VAL:HG21	1:C:99:LEU:HD13	1.98	0.46
1:D:220:THR:CG2	1:D:223:MET:HB2	2.46	0.46
1:A:153:ASN:O	1:A:156:ALA:HB3	2.16	0.46
1:D:161:VAL:HG23	1:D:275:TRP:CZ2	2.50	0.46
1:B:66:GLY:O	1:B:67:TYR:C	2.54	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:24:GLU:HB3	1:B:82:ASP:CG	2.37	0.45
1:B:109:LEU:HD12	1:B:109:LEU:N	2.31	0.45
1:A:262:ARG:O	1:A:266:VAL:HG23	2.15	0.45
1:A:29:ALA:HA	1:A:50:ASP:HB3	1.99	0.45
1:D:248:ILE:HD12	1:D:248:ILE:N	2.32	0.45
1:A:29:ALA:HB3	1:A:86:THR:HB	1.98	0.45
1:C:164:MET:SD	1:C:209:LEU:HD11	2.56	0.45
1:D:164:MET:HG3	1:D:176:VAL:HG13	1.99	0.45
1:D:216:VAL:HG11	1:D:249:ILE:HD12	1.97	0.45
1:D:207:GLU:CD	1:D:207:GLU:H	2.20	0.45
1:B:269:SER:O	1:B:273:GLN:HG3	2.16	0.45
1:D:41:GLY:HA2	2:D:296:HOH:O	2.17	0.45
1:A:224:VAL:O	1:A:225:ASP:C	2.55	0.45
1:C:256:ILE:HD12	1:C:256:ILE:C	2.37	0.45
1:D:220:THR:HG23	1:D:223:MET:CB	2.48	0.45
1:B:145:GLN:NE2	1:C:88:GLN:NE2	2.65	0.44
1:D:221:SER:HB3	1:D:250:THR:CG2	2.47	0.44
1:C:86:THR:O	1:C:108:THR:HA	2.15	0.44
1:A:178:GLY:N	1:A:200:GLN:O	2.45	0.44
1:B:198:HIS:HD2	1:B:204:TYR:OH	2.00	0.44
1:A:255:LEU:HB3	1:A:266:VAL:HG21	1.99	0.44
1:B:24:GLU:HB3	1:B:82:ASP:OD2	2.18	0.44
1:B:70:GLN:CA	1:B:92:PRO:HD3	2.47	0.44
1:D:86:THR:O	1:D:108:THR:HA	2.18	0.44
1:A:68:TRP:HE3	1:A:90:ALA:HA	1.83	0.44
1:B:25:ARG:NH1	1:B:47:VAL:HG11	2.33	0.43
1:B:83:SER:HA	1:B:105:ASN:O	2.17	0.43
1:B:183:ALA:HB1	1:B:256:ILE:HD11	2.00	0.43
1:C:216:VAL:HG11	1:C:249:ILE:HD12	2.00	0.43
1:A:68:TRP:HA	1:A:71:LEU:HD13	2.00	0.43
1:B:100:ARG:NH1	1:B:106:VAL:HB	2.33	0.43
1:B:25:ARG:HB2	1:B:82:ASP:OD1	2.19	0.43
1:C:277:GLN:NE2	1:C:277:GLN:HA	2.34	0.43
1:D:55:TYR:CG	1:D:56:PRO:HA	2.53	0.43
1:D:243:TRP:CA	1:D:248:ILE:HD11	2.44	0.43
1:A:62:LEU:O	1:A:64:HIS:CD2	2.70	0.43
1:B:32:SER:O	1:B:36:LEU:HG	2.18	0.43
1:B:80:ARG:N	1:B:81:PRO:CD	2.81	0.43
1:C:251:VAL:HG22	1:C:252:ASP:N	2.34	0.43
1:A:165:PHE:CE2	1:A:167:LEU:HB2	2.54	0.43
1:B:40:MET:O	1:B:135:GLN:HG2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:55:TYR:CG	1:B:56:PRO:HA	2.54	0.43
1:B:151:GLN:O	1:B:154:VAL:HB	2.19	0.43
1:D:165:PHE:CE2	1:D:167:LEU:HB2	2.54	0.43
1:B:27:VAL:HB	1:B:84:VAL:HG22	2.01	0.42
1:B:67:TYR:O	1:B:70:GLN:N	2.42	0.42
1:B:23:ALA:N	1:B:83:SER:OG	2.47	0.42
1:C:62:LEU:O	1:C:64:HIS:HD2	2.03	0.42
1:A:146:ARG:NH1	2:A:323:HOH:O	2.52	0.42
1:B:217:ILE:HD12	1:B:240:THR:HG21	2.01	0.42
1:A:171:GLY:O	1:C:223:MET:HB2	2.19	0.42
1:B:71:LEU:N	1:B:91:GLY:O	2.53	0.42
1:B:171:GLY:O	1:D:168:SER:HB3	2.19	0.42
1:B:92:PRO:HB2	1:B:95:VAL:CG2	2.48	0.42
1:A:112:VAL:HG22	1:A:113:PRO:N	2.33	0.42
1:B:71:LEU:CD1	1:B:76:ILE:HD11	2.47	0.42
1:C:270:LEU:O	1:C:274:LEU:HG	2.19	0.42
1:B:65:ILE:HD12	1:B:76:ILE:HG12	2.00	0.42
1:B:86:THR:OG1	1:B:87:TRP:N	2.51	0.42
1:B:94:ILE:HG22	1:B:98:GLN:HG3	2.02	0.42
1:C:144:ASN:O	1:C:148:GLU:HB2	2.20	0.42
1:C:55:TYR:CE1	1:C:57:PRO:HD3	2.54	0.42
1:C:164:MET:SD	1:C:209:LEU:CD1	3.08	0.42
1:B:230:ARG:NH1	2:B:325:HOH:O	2.53	0.41
1:C:111:ARG:HG3	1:C:111:ARG:HH11	1.85	0.41
1:B:147:LEU:O	1:B:151:GLN:HG2	2.18	0.41
1:B:224:VAL:O	1:B:225:ASP:C	2.57	0.41
1:A:49:VAL:O	1:A:64:HIS:HA	2.20	0.41
1:C:227:ASP:OD2	1:C:229:ASN:HB2	2.20	0.41
1:D:126:LEU:HD22	1:D:130:LEU:HD11	2.01	0.41
1:A:109:LEU:HD12	1:A:109:LEU:N	2.35	0.41
1:A:24:GLU:HB3	1:A:82:ASP:OD2	2.20	0.41
1:C:111:ARG:HG3	1:C:111:ARG:NH1	2.35	0.41
1:B:178:GLY:N	1:B:200:GLN:O	2.50	0.41
1:A:74:GLU:HG3	1:C:61:LYS:HA	2.02	0.41
1:B:159:ALA:HA	1:B:160:PRO:HD3	1.84	0.41
1:D:68:TRP:CZ3	1:D:69:LYS:HG3	2.56	0.41
1:B:90:ALA:HB1	1:B:96:LEU:HD21	2.03	0.41
1:B:183:ALA:CB	1:B:256:ILE:HD11	2.51	0.41
1:A:23:ALA:N	1:A:83:SER:HG	2.19	0.40
1:B:51:GLU:HB2	2:B:303:HOH:O	2.20	0.40
1:B:240:THR:HG23	1:B:242:ALA:H	1.85	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:67:TYR:HB2	1:D:70:GLN:CG	2.41	0.40
1:B:232:ARG:HH11	1:B:232:ARG:HG2	1.87	0.40
1:D:221:SER:OG	1:D:222:GLN:NE2	2.53	0.40
1:D:264:ALA:HB3	2:D:316:HOH:O	2.22	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	253/256 (99%)	238 (94%)	12 (5%)	3 (1%)	13	5
1	B	253/256 (99%)	231 (91%)	19 (8%)	3 (1%)	13	5
1	C	253/256 (99%)	242 (96%)	11 (4%)	0	100	100
1	D	253/256 (99%)	237 (94%)	15 (6%)	1 (0%)	34	24
All	All	1012/1024 (99%)	948 (94%)	57 (6%)	7 (1%)	22	12

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	91	GLY
1	B	90	ALA
1	D	200	GLN
1	A	225	ASP
1	B	67	TYR
1	A	226	GLY
1	B	226	GLY

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	207/207 (100%)	201 (97%)	6 (3%)	42	35
1	B	207/207 (100%)	200 (97%)	7 (3%)	37	30
1	C	207/207 (100%)	204 (99%)	3 (1%)	67	65
1	D	207/207 (100%)	200 (97%)	7 (3%)	37	30
All	All	828/828 (100%)	805 (97%)	23 (3%)	43	37

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	70	GLN
1	A	105	ASN
1	A	112	VAL
1	A	130	LEU
1	A	151	GLN
1	A	194	ASN
1	B	44	GLU
1	B	105	ASN
1	B	112	VAL
1	B	135	GLN
1	B	194	ASN
1	B	237	ILE
1	B	240	THR
1	C	94	ILE
1	C	194	ASN
1	C	238	THR
1	D	73	SER
1	D	94	ILE
1	D	126	LEU
1	D	153	ASN
1	D	194	ASN
1	D	238	THR
1	D	277	GLN

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

Mol	Chain	Res	Type
1	A	64	HIS
1	A	105	ASN
1	A	142	GLN
1	A	151	GLN
1	A	194	ASN
1	A	198	HIS
1	A	254	ASN
1	A	272	GLN
1	B	64	HIS
1	B	98	GLN
1	B	105	ASN
1	B	131	GLN
1	B	145	GLN
1	B	194	ASN
1	B	198	HIS
1	B	229	ASN
1	B	254	ASN
1	B	271	HIS
1	B	272	GLN
1	C	64	HIS
1	C	98	GLN
1	C	105	ASN
1	C	194	ASN
1	C	245	ASN
1	C	254	ASN
1	C	271	HIS
1	C	277	GLN
1	D	98	GLN
1	D	105	ASN
1	D	153	ASN
1	D	194	ASN
1	D	198	HIS
1	D	229	ASN
1	D	245	ASN
1	D	271	HIS
1	D	277	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

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	255/256 (99%)	0.45	15 (5%) 22 24	28, 42, 62, 80	0
1	B	255/256 (99%)	0.87	40 (15%) 2 1	28, 48, 97, 104	0
1	C	255/256 (99%)	0.38	16 (6%) 20 21	31, 47, 69, 88	0
1	D	255/256 (99%)	0.28	11 (4%) 35 38	27, 42, 62, 87	0
All	All	1020/1024 (99%)	0.49	82 (8%) 12 13	27, 45, 79, 104	0

All (82) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	96	LEU	7.4
1	A	159	ALA	7.3
1	C	156	ALA	6.8
1	B	159	ALA	6.8
1	C	277	GLN	6.4
1	C	23	ALA	6.4
1	B	92	PRO	6.1
1	B	104	VAL	5.8
1	B	23	ALA	5.6
1	B	101	ALA	5.6
1	A	155	ALA	5.5
1	A	93	GLN	5.3
1	B	67	TYR	5.3
1	D	277	GLN	5.3
1	B	99	LEU	5.2
1	B	94	ILE	5.1
1	B	93	GLN	5.0
1	B	71	LEU	4.9
1	B	102	GLN	4.9
1	A	154	VAL	4.7
1	B	130	LEU	4.6

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Mol	Chain	Res	Type	RSRZ
1	D	159	ALA	4.5
1	B	77	LEU	4.4
1	B	95	VAL	4.4
1	D	156	ALA	4.3
1	B	81	PRO	4.3
1	A	91	GLY	4.2
1	B	91	GLY	4.2
1	D	160	PRO	4.2
1	A	158	LYS	4.2
1	D	155	ALA	4.2
1	B	98	GLN	4.1
1	D	158	LYS	4.1
1	B	156	ALA	4.0
1	B	158	LYS	3.9
1	B	79	LEU	3.9
1	B	73	SER	3.8
1	B	72	SER	3.7
1	C	154	VAL	3.7
1	B	24	GLU	3.6
1	A	94	ILE	3.6
1	B	80	ARG	3.5
1	B	155	ALA	3.5
1	B	76	ILE	3.4
1	C	276	PRO	3.3
1	C	158	LYS	3.2
1	A	112	VAL	3.2
1	B	78	SER	3.2
1	D	276	PRO	3.1
1	B	74	GLU	3.1
1	C	155	ALA	3.0
1	C	150	VAL	3.0
1	B	128	LYS	3.0
1	C	153	ASN	2.9
1	C	272	GLN	2.9
1	C	151	GLN	2.9
1	B	90	ALA	2.9
1	A	71	LEU	2.9
1	C	152	GLN	2.8
1	A	156	ALA	2.8
1	C	157	LYS	2.8
1	A	157	LYS	2.7
1	A	96	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	160	PRO	2.6
1	A	23	ALA	2.6
1	B	25	ARG	2.5
1	C	159	ALA	2.5
1	B	70	GLN	2.4
1	B	131	GLN	2.4
1	D	23	ALA	2.4
1	B	157	LYS	2.3
1	A	169	ALA	2.2
1	B	256	ILE	2.2
1	B	75	GLY	2.2
1	B	47	VAL	2.2
1	D	220	THR	2.2
1	C	274	LEU	2.2
1	D	39	ALA	2.1
1	A	277	GLN	2.0
1	B	105	ASN	2.0
1	D	157	LYS	2.0
1	B	82	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 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.