



# Full wwPDB X-ray Structure Validation Report

Mar 1, 2014 – 02:44 AM GMT

PDB ID : 2BHM  
Title : CRYSTAL STRUCTURE OF VIRB8 FROM BRUCELLA SUIS  
Authors : Bayliss, R.; Baron, C.; Waksman, G.  
Deposited on : 2005-01-14  
Resolution : 2.40 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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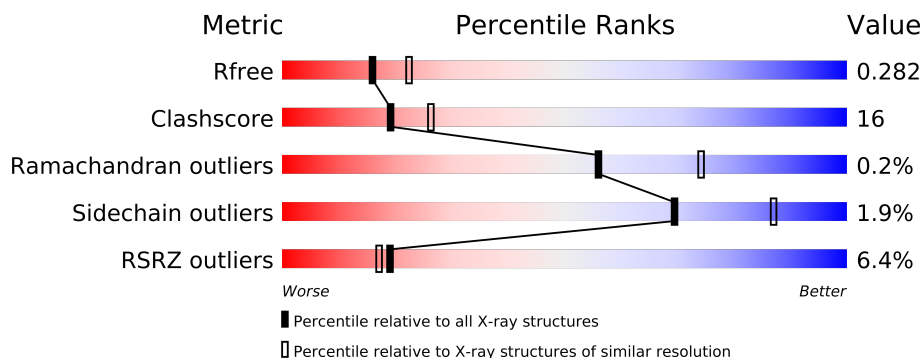
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance






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))
$R_{free}$	66092	2207 (2.40-2.40)
Clashscore	79885	2789 (2.40-2.40)
Ramachandran outliers	78287	2736 (2.40-2.40)
Sidechain outliers	78261	2737 (2.40-2.40)
RSRZ outliers	66119	2210 (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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	164	
1	B	164	
1	C	164	
1	D	164	
1	E	164	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5382 atoms, of which 0 are hydrogen and 0 are deuterium.

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 TYPE IV SECRETION SYSTEM PROTEIN VIRB8.

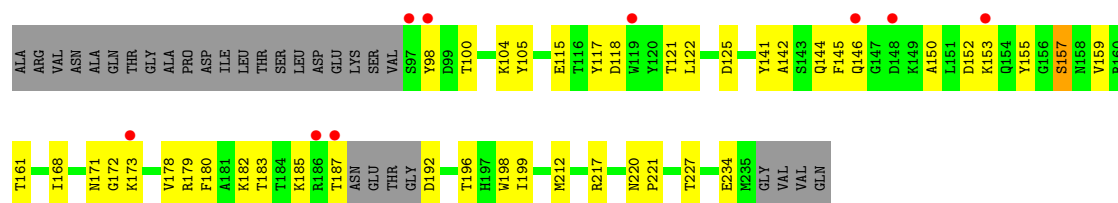
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	135	Total	C	N	O	S	0	2	1
			1077	678	179	216	4			
1	B	135	Total	C	N	O	S	0	0	1
			1063	670	176	214	3			
1	C	135	Total	C	N	O	S	0	0	1
			1063	670	176	214	3			
1	D	134	Total	C	N	O	S	0	0	1
			1055	666	175	211	3			
1	E	135	Total	C	N	O	S	0	0	1
			1063	670	176	214	3			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	16	Total	O	0	0
			16	16		
2	B	13	Total	O	0	0
			13	13		
2	C	3	Total	O	0	0
			3	3		
2	D	25	Total	O	0	0
			25	25		
2	E	4	Total	O	0	0
			4	4		



Chain E:



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	202.76Å 202.76Å 103.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	26.60 – 2.40 26.62 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.4 (26.60-2.40) 98.4 (26.62-2.40)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.71 (at 2.39Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.246 , 0.273 0.247 , 0.282	Depositor DCC
$R_{free}$ test set	2083 reflections (5.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	40.0	Xtriage
Anisotropy	0.371	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 39.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 41354 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	5382	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.78% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

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	0.47	0/1101	0.70	0/1496
1	B	0.44	0/1087	0.66	0/1479
1	C	0.41	0/1087	0.63	0/1479
1	D	0.48	0/1079	0.71	0/1468
1	E	0.41	0/1087	0.61	0/1479
All	All	0.44	0/5441	0.66	0/7401

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Close contacts

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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1077	0	1023	33	0
1	B	1063	0	1012	21	0
1	C	1063	0	1012	51	0
1	D	1055	0	1008	40	0
1	E	1063	0	1012	36	0
2	A	16	0	0	0	0
2	B	13	0	0	0	0
2	C	3	0	0	0	0
2	D	25	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	4	0	0	0	0
All	All	5382	0	5067	170	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 16.

All (170) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:150:ALA:HB3	1:C:153:LYS:HG3	1.44	0.99
1:A:196:THR:HG23	1:A:198:TRP:HE1	1.31	0.93
1:A:230:ARG:HH11	1:A:230:ARG:HG2	1.36	0.90
1:C:148:ASP:HA	1:C:153:LYS:HD2	1.53	0.90
1:C:198:TRP:CZ3	1:C:233:PRO:HG3	2.08	0.89
1:A:179:ARG:NH1	1:A:234:GLU:OE2	2.08	0.87
1:C:131:MET:SD	1:C:219:THR:HG21	2.14	0.86
1:E:199:ILE:HD11	1:E:234:GLU:HG3	1.61	0.83
1:C:195:THR:HB	1:D:179:ARG:NH1	1.95	0.82
1:C:141:TYR:O	1:C:144:GLN:HG2	1.83	0.79
1:A:139[A]:GLN:HG3	1:A:140:SER:N	1.99	0.77
1:C:222:LEU:HD21	1:E:98:TYR:HE1	1.49	0.76
1:B:199:ILE:HD11	1:B:234:GLU:HG3	1.66	0.76
1:A:230:ARG:HG2	1:A:230:ARG:NH1	2.00	0.75
1:C:222:LEU:HD21	1:E:98:TYR:CE1	2.21	0.75
1:A:230:ARG:NH1	1:A:232:ASP:OD2	2.22	0.73
1:E:196:THR:HB	1:E:198:TRP:HE1	1.56	0.70
1:A:196:THR:CG2	1:A:198:TRP:HE1	2.03	0.70
1:C:196:THR:HB	1:C:198:TRP:HE1	1.57	0.69
1:C:222:LEU:CD2	1:E:98:TYR:HE1	2.06	0.68
1:C:118:ASP:HB3	1:C:121:THR:HB	1.75	0.67
1:B:196:THR:HB	1:B:198:TRP:HE1	1.60	0.65
1:C:171:ASN:ND2	1:C:175:ILE:HB	2.11	0.65
1:C:198:TRP:HZ3	1:C:233:PRO:HG3	1.62	0.65
1:A:171:ASN:ND2	1:A:175:ILE:HB	2.11	0.65
1:E:171:ASN:HD21	1:E:173:LYS:HB2	1.61	0.65
1:A:139[A]:GLN:NE2	1:D:175:ILE:HD11	2.12	0.64
1:D:196:THR:HB	1:D:198:TRP:HE1	1.61	0.64
1:C:171:ASN:HD21	1:C:175:ILE:HB	1.62	0.63
1:C:199:ILE:HD11	1:C:234:GLU:CD	2.19	0.62
1:D:230:ARG:HG2	1:D:230:ARG:HH11	1.65	0.61
1:B:196:THR:HG21	1:B:198:TRP:CZ2	2.34	0.61
1:D:212:MET:HG3	1:D:217:ARG:HG3	1.81	0.61
1:D:179:ARG:NH1	1:D:234:GLU:OE2	2.34	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:115:GLU:HB3	1:B:182:LYS:HE3	1.83	0.61
1:D:117:TYR:OH	1:D:152:ASP:HA	2.01	0.60
1:A:230:ARG:NH2	1:D:144:GLN:OE1	2.34	0.60
1:A:111:VAL:HG21	1:A:178:VAL:HG11	1.82	0.60
1:B:158:ASN:HA	1:B:187:THR:HB	1.84	0.59
1:D:199:ILE:HD11	1:D:234:GLU:OE2	2.02	0.59
1:C:199:ILE:HD11	1:C:234:GLU:OE2	2.02	0.59
1:E:142:ALA:O	1:E:146:GLN:HG3	2.02	0.59
1:C:155:TYR:HD2	1:C:159:VAL:HG21	1.68	0.59
1:C:155:TYR:HD2	1:C:159:VAL:CG2	2.15	0.59
1:E:196:THR:HB	1:E:198:TRP:NE1	2.17	0.59
1:D:230:ARG:NH1	1:D:232:ASP:OD1	2.36	0.58
1:E:196:THR:HG21	1:E:198:TRP:CZ2	2.39	0.57
1:A:230:ARG:HH12	1:A:232:ASP:CG	2.07	0.57
1:B:212:MET:HG3	1:B:217:ARG:HG3	1.88	0.56
1:A:121:THR:HG22	1:A:125:ASP:OD2	2.05	0.56
1:C:121:THR:HG22	1:C:125:ASP:OD2	2.06	0.56
1:C:155:TYR:HD2	1:C:159:VAL:CB	2.19	0.56
1:A:209:PRO:HA	1:A:212:MET:CE	2.35	0.56
1:E:141:TYR:O	1:E:144:GLN:HG2	2.06	0.56
1:C:155:TYR:HD2	1:C:159:VAL:HB	1.72	0.55
1:D:137:GLU:OE2	1:D:228:SER:HA	2.07	0.54
1:D:230:ARG:CG	1:D:230:ARG:HH11	2.20	0.54
1:C:155:TYR:CD2	1:C:159:VAL:HB	2.41	0.54
1:D:171:ASN:C	1:D:171:ASN:OD1	2.47	0.54
1:E:185:LYS:HE2	1:E:192:ASP:HA	1.89	0.53
1:B:171:ASN:ND2	1:B:175:ILE:HB	2.23	0.53
1:E:212:MET:HG3	1:E:217:ARG:HG3	1.89	0.53
1:C:195:THR:HB	1:D:179:ARG:HH12	1.74	0.53
1:D:200:ALA:HA	1:D:230:ARG:O	2.08	0.53
1:D:209:PRO:HA	1:D:212:MET:CE	2.39	0.53
1:A:139[A]:GLN:HE22	1:D:175:ILE:HD11	1.73	0.53
1:E:179:ARG:NH1	1:E:234:GLU:OE2	2.41	0.52
1:A:209:PRO:HA	1:A:212:MET:HE2	1.91	0.52
1:B:179:ARG:HD3	1:B:234:GLU:OE2	2.09	0.52
1:E:100:THR:HG22	1:E:104:LYS:HE3	1.91	0.52
1:C:148:ASP:HA	1:C:153:LYS:CD	2.35	0.52
1:A:230:ARG:NH1	1:A:230:ARG:CG	2.69	0.51
1:C:117:TYR:OH	1:C:152:ASP:HA	2.09	0.51
1:E:122:LEU:HD11	1:E:145:PHE:CE2	2.45	0.51
1:A:161:THR:HA	1:A:183:THR:O	2.09	0.51
1:D:118:ASP:HB3	1:D:121:THR:HB	1.93	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:209:PRO:O	1:B:212:MET:HG2	2.10	0.51
1:D:171:ASN:OD1	1:D:172:GLY:N	2.43	0.51
1:E:157:SER:O	1:E:187:THR:OG1	2.28	0.51
1:C:121:THR:HG22	1:C:125:ASP:CG	2.32	0.50
1:D:209:PRO:O	1:D:212:MET:HG2	2.11	0.50
1:A:200:ALA:HA	1:A:230:ARG:O	2.11	0.49
1:D:141:TYR:O	1:D:144:GLN:HG2	2.12	0.49
1:E:117:TYR:OH	1:E:152:ASP:HA	2.12	0.49
1:C:196:THR:HB	1:C:198:TRP:NE1	2.25	0.49
1:A:119:TRP:HZ3	1:A:152:ASP:HB2	1.78	0.49
1:C:200:ALA:HB2	1:C:231:VAL:HG22	1.93	0.49
1:C:102:MET:HG2	1:C:222:LEU:CD1	2.43	0.49
1:C:200:ALA:HA	1:C:230:ARG:O	2.12	0.49
1:A:212:MET:HG3	1:A:217:ARG:HG3	1.94	0.49
1:C:222:LEU:CD2	1:E:98:TYR:CE1	2.88	0.48
1:C:168:ILE:HG12	1:C:178:VAL:HG22	1.95	0.48
1:D:161:THR:HA	1:D:183:THR:O	2.13	0.48
1:D:209:PRO:HA	1:D:212:MET:HE2	1.94	0.48
1:A:196:THR:HG23	1:A:198:TRP:NE1	2.14	0.48
1:B:196:THR:HB	1:B:198:TRP:NE1	2.27	0.48
1:E:121:THR:HG22	1:E:125:ASP:CG	2.34	0.48
1:D:196:THR:HG21	1:D:198:TRP:CZ2	2.49	0.48
1:E:121:THR:HG22	1:E:125:ASP:OD2	2.14	0.48
1:B:220:ASN:N	1:B:221:PRO:HD3	2.29	0.47
1:C:127:GLU:O	1:C:131:MET:HG2	2.15	0.47
1:D:179:ARG:HG3	2:D:2015:HOH:O	2.14	0.47
1:A:171:ASN:OD1	1:A:173:LYS:N	2.37	0.47
1:C:145:PHE:HA	1:C:150:ALA:HB1	1.97	0.47
1:D:111:VAL:HG21	1:D:178:VAL:HG11	1.97	0.47
1:D:119:TRP:HZ3	1:D:152:ASP:HB2	1.79	0.47
1:E:220:ASN:N	1:E:221:PRO:HD3	2.31	0.46
1:D:157:SER:O	1:D:187:THR:OG1	2.32	0.46
1:D:161:THR:CG2	1:D:182:LYS:HG2	2.46	0.46
1:E:115:GLU:HB3	1:E:182:LYS:HE3	1.98	0.46
1:B:198:TRP:CZ3	1:B:233:PRO:HG3	2.50	0.46
1:C:221:PRO:HG2	1:E:105:TYR:CG	2.51	0.46
1:D:201:THR:HB	1:D:230:ARG:HB3	1.99	0.45
1:C:148:ASP:CA	1:C:153:LYS:HD2	2.36	0.45
1:B:142:ALA:O	1:B:146:GLN:HG3	2.16	0.45
1:B:171:ASN:OD1	1:B:173:LYS:N	2.50	0.45
1:C:120:TYR:N	1:C:120:TYR:CD1	2.84	0.45
1:C:118:ASP:O	1:C:119:TRP:C	2.55	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:144:GLN:CD	1:D:230:ARG:NH2	2.70	0.45
1:E:100:THR:O	1:E:104:LYS:HG3	2.17	0.45
1:A:230:ARG:NH1	1:A:232:ASP:CG	2.68	0.45
1:B:207:VAL:O	1:B:208:ASN:C	2.55	0.45
1:E:168:ILE:HG12	1:E:178:VAL:HG22	1.99	0.44
1:E:155:TYR:HB3	1:E:159:VAL:HB	1.98	0.44
1:B:151:LEU:HD22	1:B:155:TYR:CE1	2.52	0.44
1:C:196:THR:HG21	1:C:198:TRP:CZ2	2.53	0.44
1:D:115:GLU:HG3	1:D:180:PHE:CZ	2.53	0.44
1:D:131:MET:HE2	1:D:131:MET:HB2	1.78	0.44
1:B:115:GLU:HG3	1:B:180:PHE:CE1	2.53	0.44
1:D:182:LYS:O	1:D:195:THR:HA	2.18	0.44
1:C:125:ASP:O	1:C:129:VAL:HG23	2.16	0.44
1:C:151:LEU:O	1:C:155:TYR:HD1	2.01	0.43
1:C:108:SER:O	1:C:112:ILE:HG13	2.17	0.43
1:B:172:GLY:O	1:B:173:LYS:HG3	2.18	0.43
1:E:161:THR:HA	1:E:183:THR:O	2.19	0.43
1:E:171:ASN:CG	1:E:172:GLY:N	2.71	0.43
1:C:155:TYR:CD2	1:C:159:VAL:HG21	2.49	0.43
1:A:115:GLU:HG3	1:A:180:PHE:CZ	2.54	0.43
1:E:161:THR:CG2	1:E:182:LYS:HG2	2.48	0.43
1:D:104:LYS:HE2	1:D:168:ILE:O	2.19	0.43
1:A:133:SER:HB3	1:A:226:VAL:HB	2.00	0.42
1:D:218:LEU:HD23	1:D:218:LEU:HA	1.79	0.42
1:C:217:ARG:O	1:C:221:PRO:HG3	2.19	0.42
1:B:198:TRP:HZ3	1:B:233:PRO:HG3	1.83	0.42
1:C:116:THR:HG22	1:C:117:TYR:N	2.34	0.42
1:E:122:LEU:HD11	1:E:145:PHE:CZ	2.53	0.42
1:C:179:ARG:NH1	1:C:234:GLU:OE2	2.51	0.42
1:C:220:ASN:N	1:C:221:PRO:HD3	2.35	0.42
1:E:155:TYR:CD2	1:E:155:TYR:N	2.84	0.42
1:C:100:THR:O	1:C:104:LYS:HG3	2.19	0.42
1:D:230:ARG:NH1	1:D:230:ARG:CG	2.81	0.42
1:B:217:ARG:O	1:B:221:PRO:HG3	2.20	0.42
1:A:117:TYR:OH	1:A:152:ASP:HA	2.18	0.42
1:C:232:ASP:HA	1:C:233:PRO:HD3	1.84	0.42
1:E:150:ALA:O	1:E:153:LYS:HB2	2.19	0.42
1:D:220:ASN:N	1:D:221:PRO:HD3	2.34	0.42
1:E:115:GLU:HG3	1:E:180:PHE:CZ	2.55	0.41
1:E:199:ILE:HD11	1:E:234:GLU:CG	2.41	0.41
1:A:209:PRO:O	1:A:212:MET:HG2	2.21	0.41
1:C:105:TYR:O	1:C:109:GLN:HG2	2.21	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:171:ASN:HD21	1:A:175:ILE:HB	1.85	0.41
1:E:118:ASP:HB3	1:E:121:THR:HB	2.02	0.41
1:E:171:ASN:OD1	1:E:172:GLY:N	2.50	0.41
1:D:121:THR:O	1:D:121:THR:HG22	2.20	0.41
1:D:131:MET:HE1	1:D:219:THR:HB	2.03	0.41
1:A:182:LYS:O	1:A:195:THR:HA	2.21	0.41
1:A:137:GLU:OE2	1:A:228:SER:HA	2.20	0.41
1:C:121:THR:HG22	1:C:121:THR:O	2.22	0.40
1:B:209:PRO:HA	1:B:212:MET:CE	2.52	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

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	133/164 (81%)	130 (98%)	3 (2%)	0	100	100
1	B	131/164 (80%)	127 (97%)	4 (3%)	0	100	100
1	C	131/164 (80%)	124 (95%)	6 (5%)	1 (1%)	27	39
1	D	130/164 (79%)	126 (97%)	4 (3%)	0	100	100
1	E	131/164 (80%)	127 (97%)	4 (3%)	0	100	100
All	All	656/820 (80%)	634 (97%)	21 (3%)	1 (0%)	56	74

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	172	GLY

### 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	118/142 (83%)	115 (98%)	3 (2%)	60	80
1	B	117/142 (82%)	116 (99%)	1 (1%)	87	96
1	C	117/142 (82%)	115 (98%)	2 (2%)	73	89
1	D	116/142 (82%)	113 (97%)	3 (3%)	59	79
1	E	117/142 (82%)	115 (98%)	2 (2%)	73	89
All	All	585/710 (82%)	574 (98%)	11 (2%)	69	87

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	99	ASP
1	A	157	SER
1	A	212	MET
1	B	154	GLN
1	C	173	LYS
1	C	187	THR
1	D	157	SER
1	D	173	LYS
1	D	194	GLU
1	E	157	SER
1	E	227	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	205	GLN
1	A	225	ASN
1	C	146	GLN
1	E	205	GLN
1	E	225	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

There are no ligands in this entry.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	135/164 (82%)	-0.01	4 (2%)	48	45	20, 30, 67, 79	2 (1%)
1	B	135/164 (82%)	0.04	6 (4%)	33	30	23, 38, 66, 86	0
1	C	135/164 (82%)	0.67	20 (14%)	3	2	32, 59, 89, 94	0
1	D	134/164 (81%)	-0.07	4 (2%)	48	45	20, 33, 65, 81	0
1	E	135/164 (82%)	0.39	9 (6%)	17	16	31, 53, 83, 93	0
All	All	674/820 (82%)	0.20	43 (6%)	19	17	20, 45, 80, 94	2 (0%)

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	97	SER	7.1
1	E	187	THR	6.3
1	C	98	TYR	6.2
1	C	147	GLY	5.4
1	C	187	THR	4.7
1	C	148	ASP	4.5
1	E	98	TYR	4.4
1	C	185	LYS	4.3
1	B	187	THR	4.2
1	B	186	ARG	3.8
1	C	153	LYS	3.7
1	C	150	ALA	3.7
1	D	187	THR	3.6
1	E	186	ARG	3.4
1	E	173	LYS	3.4
1	A	148	ASP	3.1
1	C	119	TRP	3.1
1	C	172	GLY	3.0
1	B	192	ASP	3.0
1	D	148	ASP	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	154	GLN	2.8
1	E	119	TRP	2.8
1	B	173	LYS	2.8
1	A	186	ARG	2.8
1	C	152	ASP	2.8
1	C	97	SER	2.7
1	C	139	GLN	2.7
1	C	146	GLN	2.7
1	A	155	TYR	2.7
1	B	98	TYR	2.7
1	C	192	ASP	2.6
1	C	155	TYR	2.6
1	B	185	LYS	2.6
1	C	149	LYS	2.6
1	C	131	MET	2.5
1	D	186	ARG	2.4
1	C	173	LYS	2.4
1	C	186	ARG	2.3
1	A	119	TRP	2.2
1	E	148	ASP	2.2
1	E	153	LYS	2.1
1	E	146	GLN	2.0
1	C	120	TYR	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

There are no ligands in this entry.

## 6.5 Other polymers ⓘ

There are no such residues in this entry.