



Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 07:15 pm BST

PDB ID : 6V78
Title : OmpK37 porin
Authors : Rocker, A.; Lithgow, T.
Deposited on : 2019-12-08
Resolution : 2.60 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 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 : 2.11
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) : 2.11

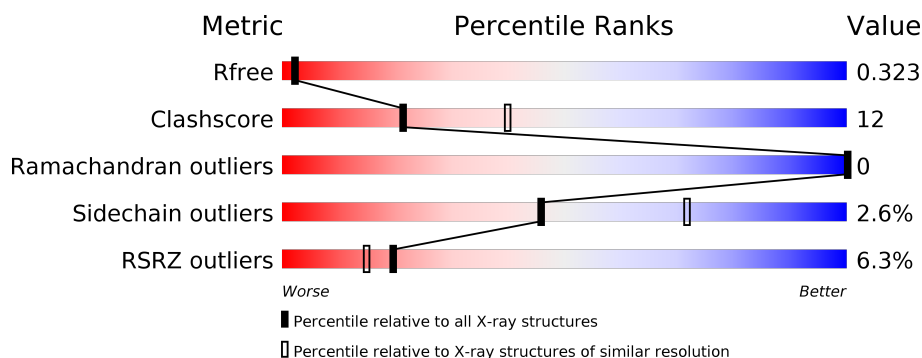
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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 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	356	<div> <div>4%</div> <div>71%</div> <div>28%</div> <div>..</div> </div>
1	B	356	<div> <div>11%</div> <div>69%</div> <div>30%</div> <div>.</div> </div>
1	C	356	<div> <div>3%</div> <div>76%</div> <div>22%</div> <div>..</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8455 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 OmpK37.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	353	Total	C	N	O	S	0	0	0
			2794	1746	463	576	9			
1	B	353	Total	C	N	O	S	0	0	0
			2794	1746	463	576	9			
1	C	353	Total	C	N	O	S	0	0	0
			2794	1746	463	576	9			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP S5UCA2
A	-1	ALA	-	expression tag	UNP S5UCA2
A	0	MET	-	expression tag	UNP S5UCA2
B	-2	GLY	-	expression tag	UNP S5UCA2
B	-1	ALA	-	expression tag	UNP S5UCA2
B	0	MET	-	expression tag	UNP S5UCA2
C	-2	GLY	-	expression tag	UNP S5UCA2
C	-1	ALA	-	expression tag	UNP S5UCA2
C	0	MET	-	expression tag	UNP S5UCA2

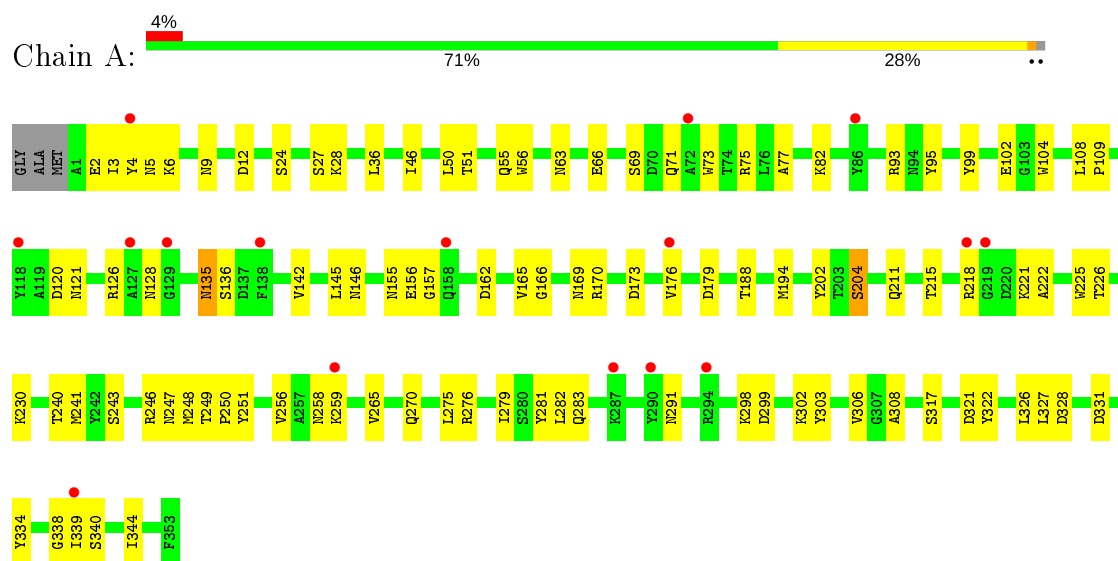
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	25	Total	O	0	0
			25	25		
2	B	15	Total	O	0	0
			15	15		
2	C	33	Total	O	0	0
			33	33		

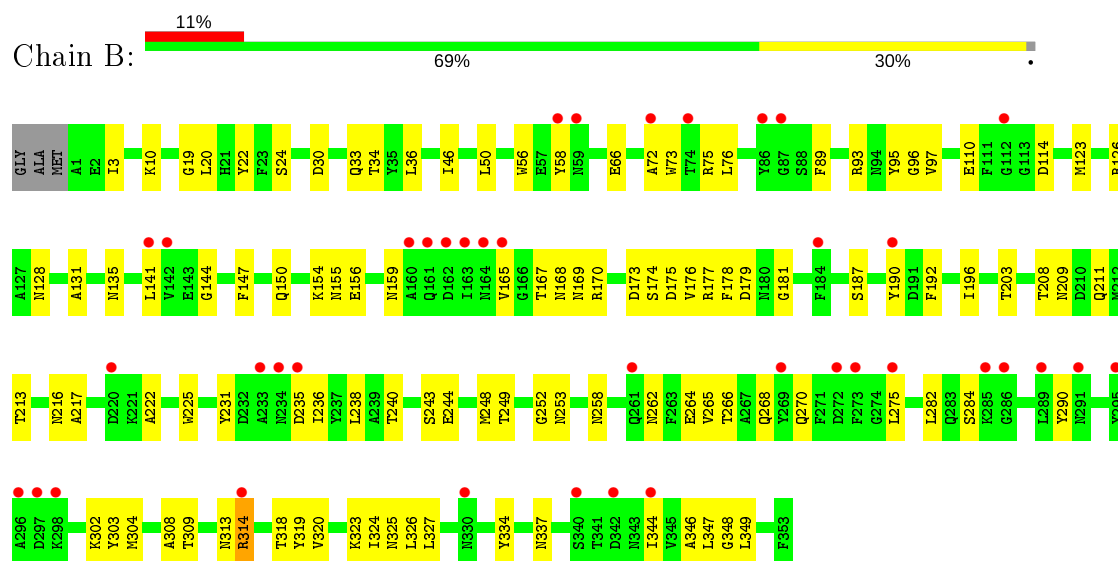
3 Residue-property plots [i](#)

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.

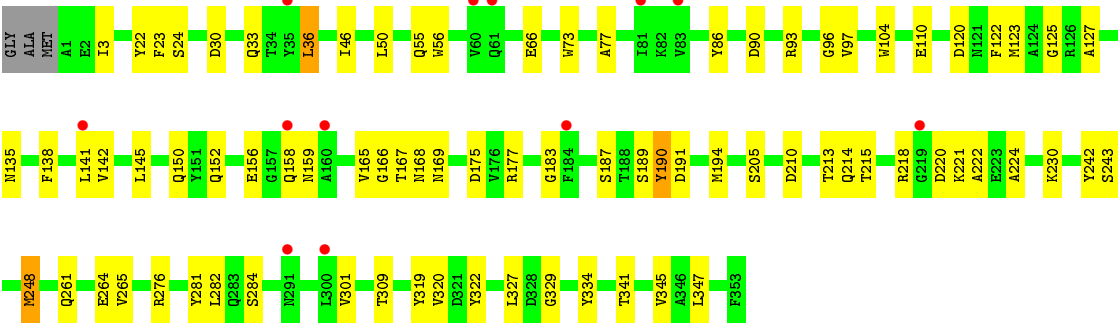
• Molecule 1: OmpK37



• Molecule 1: OmpK37



• Molecule 1: OmpK37



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	109.52Å 138.51Å 91.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.34 – 2.60 49.34 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.3 (49.34-2.60) 99.2 (49.34-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.69 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.14 _3260	Depositor
R, R_{free}	0.257 , 0.323 0.257 , 0.323	Depositor DCC
R_{free} test set	2167 reflections (4.46%)	wwPDB-VP
Wilson B-factor (Å ²)	45.2	Xtriage
Anisotropy	0.760	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 41.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	8455	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.12% 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

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	0.55	0/2855	0.70	0/3860
1	B	0.42	0/2855	0.64	0/3860
1	C	0.46	0/2855	0.68	0/3860
All	All	0.48	0/8565	0.67	0/11580

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

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	2794	0	2552	67	0
1	B	2794	0	2552	79	0
1	C	2794	0	2552	60	0
2	A	25	0	0	1	0
2	B	15	0	0	0	0
2	C	33	0	0	3	0
All	All	8455	0	7656	192	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 12.

All (192) 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:C:97:VAL:HG22	1:C:187:SER:HB3	1.54	0.89
1:C:96:GLY:HA2	1:C:150:GLN:HE22	1.42	0.84
1:C:213:THR:HG23	1:C:214:GLN:HG3	1.61	0.81
1:B:154:LYS:HG3	1:B:181:GLY:HA2	1.68	0.75
1:B:309:THR:HG23	1:B:319:TYR:HB3	1.68	0.74
1:C:123:MET:HG2	1:C:150:GLN:HE21	1.57	0.70
1:B:264:GLU:HG2	1:B:282:LEU:HG	1.73	0.70
1:B:175:ASP:OD2	1:B:177:ARG:NH1	2.26	0.69
1:C:222:ALA:HB1	1:C:248:MET:HB2	1.76	0.68
1:B:222:ALA:HB1	1:B:248:MET:HB3	1.76	0.67
1:B:3:ILE:HD12	1:C:3:ILE:HG21	1.78	0.65
1:A:176:VAL:HG22	1:A:211:GLN:HG3	1.79	0.65
1:A:156:GLU:HG2	1:A:166:GLY:HA2	1.80	0.64
1:C:220:ASP:OD1	1:C:221:LYS:HG3	1.97	0.64
1:C:224:ALA:HB2	1:C:248:MET:HE1	1.78	0.63
1:C:158:GLN:NE2	1:C:167:THR:O	2.25	0.63
1:C:158:GLN:HE22	1:C:167:THR:C	2.02	0.62
1:B:196:ILE:HG13	1:B:231:TYR:HD1	1.63	0.62
1:B:24:SER:OG	1:B:344:ILE:HA	2.00	0.62
1:C:210:ASP:HA	1:C:213:THR:HG22	1.81	0.62
1:A:69:SER:O	1:A:71:GLN:NE2	2.29	0.62
1:B:97:VAL:H	1:B:150:GLN:HE22	1.47	0.61
1:C:329:GLY:HA2	1:C:341:THR:HG21	1.82	0.60
1:B:325:ASN:OD1	1:B:327:LEU:HD12	2.01	0.60
1:B:266:THR:HB	1:B:268:GLN:HE21	1.65	0.60
1:C:156:GLU:HG2	1:C:166:GLY:HA2	1.84	0.60
1:A:246:ARG:HH11	1:A:259:LYS:HE3	1.67	0.60
1:A:222:ALA:HB1	1:A:248:MET:HB3	1.84	0.59
1:A:279:ILE:HG23	1:A:306:VAL:HG22	1.85	0.59
1:B:58:TYR:CZ	1:B:72:ALA:HB1	2.37	0.59
1:C:309:THR:HG23	1:C:319:TYR:HB3	1.84	0.58
1:A:95:TYR:OH	1:A:126:ARG:NH1	2.36	0.58
1:B:159:ASN:HB2	1:B:165:VAL:HG21	1.85	0.58
1:C:158:GLN:HG3	1:C:165:VAL:O	2.04	0.58
1:A:298:LYS:NZ	1:A:331:ASP:OD2	2.25	0.58
1:A:63:ASN:O	1:C:169:ASN:HB2	2.05	0.57
1:B:170:ARG:HG2	1:B:179:ASP:OD1	2.05	0.57
1:B:97:VAL:H	1:B:150:GLN:NE2	2.03	0.56
1:B:302:LYS:HG2	1:B:326:LEU:HB2	1.87	0.56
1:A:327:LEU:HB3	1:A:334:TYR:CE2	2.41	0.56
1:A:51:THR:OG1	1:A:82:LYS:HB3	2.05	0.56
1:B:225:TRP:NE1	1:B:244:GLU:OE1	2.34	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:56:TRP:CD1	1:B:36:LEU:HD21	2.41	0.55
1:C:97:VAL:HG22	1:C:187:SER:CB	2.32	0.55
1:A:108:LEU:HB3	1:A:109:PRO:HD2	1.89	0.55
1:A:36:LEU:HD22	1:C:56:TRP:CD1	2.42	0.55
1:B:175:ASP:HB3	1:B:178:PHE:HD2	1.70	0.54
1:B:216:ASN:HB2	1:B:290:TYR:HB3	1.87	0.54
1:A:55:GLN:O	1:A:77:ALA:HA	2.08	0.54
1:B:46:ILE:HG21	1:B:50:LEU:HD23	1.89	0.54
1:C:242:TYR:OH	1:C:261:GLN:HG2	2.08	0.53
1:A:303:TYR:HA	1:A:326:LEU:HG	1.89	0.53
1:C:138:PHE:O	1:C:141:LEU:HD12	2.08	0.53
1:A:66:GLU:HB3	1:C:73:TRP:CD1	2.44	0.53
1:B:93:ARG:HH21	1:B:128:ASN:ND2	2.06	0.53
1:B:156:GLU:O	1:B:167:THR:N	2.42	0.53
1:B:73:TRP:CD1	1:C:66:GLU:HB3	2.44	0.53
1:A:240:THR:HG22	1:A:265:VAL:HB	1.90	0.53
1:B:89:PHE:CE1	1:B:131:ALA:HB1	2.45	0.52
1:B:96:GLY:HA2	1:B:150:GLN:HE22	1.74	0.52
1:C:135:ASN:OD1	1:C:138:PHE:N	2.42	0.52
1:C:265:VAL:HG12	1:C:281:TYR:HB3	1.92	0.52
1:A:102:GLU:OE1	1:A:241:MET:HE3	2.10	0.52
1:B:76:LEU:HD11	1:B:95:TYR:HE1	1.75	0.52
1:A:104:TRP:CE2	1:A:230:LYS:HE2	2.44	0.52
1:A:121:ASN:OD1	1:A:250:PRO:HD3	2.10	0.52
1:B:22:TYR:OH	1:B:110:GLU:OE2	2.24	0.52
1:B:155:ASN:OD1	1:B:169:ASN:ND2	2.38	0.52
1:A:204:SER:OG	1:A:221:LYS:NZ	2.44	0.51
1:B:240:THR:HB	1:B:265:VAL:HG23	1.91	0.51
1:C:190:TYR:HD1	1:C:191:ASP:N	2.09	0.50
1:C:24:SER:O	1:C:30:ASP:HB2	2.11	0.50
1:A:328:ASP:N	1:A:334:TYR:HE2	2.09	0.50
1:B:58:TYR:OH	1:B:72:ALA:HB1	2.12	0.50
1:B:303:TYR:HA	1:B:326:LEU:HG	1.93	0.49
1:B:266:THR:HB	1:B:268:GLN:NE2	2.26	0.49
1:B:19:GLY:HA2	1:B:34:THR:HG23	1.94	0.49
1:B:313:ASN:HB2	1:B:314:ARG:NH2	2.27	0.49
1:A:162:ASP:HB3	1:A:165:VAL:HG23	1.95	0.49
1:B:3:ILE:HD12	1:C:3:ILE:CG2	2.43	0.49
1:B:304:MET:HB3	1:B:324:ILE:CG1	2.43	0.49
1:A:73:TRP:CD1	1:B:66:GLU:HB3	2.47	0.49
1:C:110:GLU:HB2	1:C:319:TYR:CE1	2.48	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:75:ARG:O	1:B:126:ARG:HD2	2.12	0.49
1:B:217:ALA:HA	1:B:258:ASN:OD1	2.12	0.49
1:B:327:LEU:HD22	1:B:334:TYR:CE2	2.48	0.49
1:B:303:TYR:CE2	1:B:323:LYS:HE2	2.48	0.49
1:A:246:ARG:NH1	1:A:259:LYS:HE3	2.27	0.48
1:B:24:SER:HG	1:B:344:ILE:HA	1.76	0.48
1:C:93:ARG:NH1	1:C:120:ASP:OD1	2.37	0.48
1:C:284:SER:HB3	1:C:301:VAL:HB	1.94	0.48
1:B:209:ASN:O	1:B:213:THR:HG23	2.13	0.48
1:C:159:ASN:HB2	1:C:165:VAL:HG21	1.95	0.48
1:A:4:TYR:CZ	1:A:6:LYS:HB3	2.49	0.47
1:C:218:ARG:O	2:C:401:HOH:O	2.20	0.47
1:C:327:LEU:HD22	1:C:334:TYR:CE2	2.49	0.47
1:B:89:PHE:HE1	1:B:131:ALA:HB1	1.78	0.47
1:C:46:ILE:HD12	1:C:50:LEU:HG	1.95	0.47
1:C:264:GLU:HG2	1:C:282:LEU:HG	1.95	0.47
1:A:135:ASN:HD22	1:A:136:SER:N	2.13	0.47
1:A:251:TYR:HD2	1:A:339:ILE:HD11	1.80	0.47
1:C:22:TYR:OH	1:C:110:GLU:OE1	2.20	0.47
1:A:218:ARG:O	1:A:247:ASN:ND2	2.48	0.46
1:A:27:SER:HB2	1:C:166:GLY:O	2.14	0.46
1:A:2:GLU:HA	1:A:12:ASP:OD1	2.16	0.46
1:A:302:LYS:HG2	1:A:326:LEU:HB2	1.96	0.46
1:B:252:GLY:HA3	1:B:337:ASN:HB3	1.97	0.46
1:A:5:ASN:HA	1:A:9:ASN:O	2.15	0.46
1:A:75:ARG:O	1:A:126:ARG:HD2	2.16	0.46
1:B:110:GLU:HB2	1:B:319:TYR:CE1	2.51	0.46
1:C:210:ASP:HA	1:C:213:THR:CG2	2.46	0.46
1:B:114:ASP:OD1	1:B:114:ASP:N	2.48	0.45
1:A:46:ILE:HB	1:A:50:LEU:HB3	1.97	0.45
1:A:157:GLY:HA2	1:A:170:ARG:HB3	1.99	0.45
1:A:241:MET:HG3	1:A:241:MET:O	2.17	0.45
1:C:138:PHE:HB3	1:C:142:VAL:CG2	2.46	0.45
1:B:236:ILE:HG22	1:B:238:LEU:HD12	1.98	0.45
1:B:10:LYS:HB3	1:B:10:LYS:HE3	1.54	0.45
1:B:156:GLU:HB3	1:B:208:THR:HG21	1.99	0.45
1:A:328:ASP:OD2	1:A:331:ASP:HB2	2.16	0.44
1:B:275:LEU:HD11	1:B:308:ALA:HB1	1.99	0.44
1:B:304:MET:HB3	1:B:324:ILE:HG12	1.99	0.44
1:B:320:VAL:HA	1:B:346:ALA:O	2.16	0.44
1:B:176:VAL:HG12	1:B:211:GLN:HG3	1.98	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:270:GLN:OE1	1:A:276:ARG:HB2	2.18	0.44
1:A:302:LYS:HG2	1:A:326:LEU:HD12	2.00	0.44
1:C:145:LEU:HD12	1:C:189:SER:O	2.17	0.44
1:A:248:MET:HG3	1:A:249:THR:N	2.31	0.44
1:B:56:TRP:CE2	1:C:36:LEU:HD11	2.51	0.44
1:C:127:ALA:HB1	1:C:152:GLN:NE2	2.32	0.44
1:C:96:GLY:HA2	1:C:150:GLN:NE2	2.22	0.44
1:A:162:ASP:HB3	1:A:165:VAL:CG2	2.48	0.44
1:A:170:ARG:HG2	1:A:179:ASP:OD1	2.17	0.44
1:A:218:ARG:H	1:A:258:ASN:HD21	1.65	0.44
1:A:317:SER:HB3	2:A:419:HOH:O	2.17	0.44
1:B:318:THR:HB	1:B:349:LEU:HD12	1.99	0.44
1:B:123:MET:HE3	1:B:203:THR:HB	1.99	0.43
1:B:304:MET:HB3	1:B:324:ILE:HD11	1.99	0.43
1:A:93:ARG:NH2	1:A:120:ASP:OD1	2.47	0.43
1:A:215:THR:HB	1:A:256:VAL:HG21	2.00	0.43
1:B:24:SER:O	1:B:30:ASP:HB2	2.19	0.43
1:B:167:THR:HG22	1:B:168:ASN:H	1.84	0.43
1:B:175:ASP:HB3	1:B:178:PHE:CD2	2.52	0.43
1:A:99:TYR:HA	1:A:102:GLU:HB3	2.01	0.43
1:A:226:THR:HB	1:A:243:SER:HB2	2.01	0.43
1:B:320:VAL:HB	1:B:347:LEU:HD12	1.99	0.43
1:C:175:ASP:OD1	1:C:177:ARG:HD3	2.19	0.43
1:C:183:GLY:HA3	1:C:205:SER:HB2	2.00	0.43
1:C:320:VAL:HG22	1:C:347:LEU:CD1	2.49	0.43
1:C:55:GLN:O	1:C:77:ALA:HA	2.18	0.43
1:A:3:ILE:HG21	1:C:3:ILE:HD12	2.00	0.43
1:A:275:LEU:HD11	1:A:308:ALA:HB1	2.00	0.43
1:B:144:GLY:O	1:B:190:TYR:HA	2.19	0.43
1:B:190:TYR:HE1	1:B:192:PHE:CD2	2.37	0.42
1:B:248:MET:HG3	1:B:249:THR:N	2.33	0.42
1:C:215:THR:HG22	2:C:417:HOH:O	2.18	0.42
1:A:146:ASN:O	1:A:188:THR:HA	2.18	0.42
1:B:167:THR:HG22	1:B:168:ASN:N	2.34	0.42
1:A:56:TRP:CD1	1:B:36:LEU:CD2	3.02	0.42
1:A:27:SER:O	1:C:158:GLN:NE2	2.52	0.42
1:C:122:PHE:N	1:C:122:PHE:CD1	2.87	0.42
1:B:19:GLY:O	1:B:348:GLY:HA2	2.19	0.42
1:C:104:TRP:CE2	1:C:230:LYS:HE2	2.55	0.42
1:B:97:VAL:N	1:B:150:GLN:HE22	2.15	0.42
1:A:108:LEU:HB3	1:A:109:PRO:CD	2.50	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:135:ASN:HB3	1:B:147:PHE:CE2	2.55	0.42
1:A:281:TYR:C	1:A:282:LEU:HD12	2.40	0.41
1:A:24:SER:HB2	1:A:344:ILE:HA	2.01	0.41
1:A:142:VAL:HG21	1:A:145:LEU:HD22	2.01	0.41
1:B:243:SER:HG	1:B:262:ASN:HD22	1.66	0.41
1:C:97:VAL:CG2	1:C:187:SER:HB3	2.38	0.41
1:A:36:LEU:CD2	1:C:56:TRP:CD1	3.03	0.41
1:B:20:LEU:HG	1:B:33:GLN:HB2	2.01	0.41
1:C:221:LYS:HE3	1:C:221:LYS:HB3	1.67	0.41
1:A:283:GLN:NE2	1:A:299:ASP:OD1	2.49	0.41
1:A:135:ASN:HD22	1:A:136:SER:H	1.67	0.41
1:A:202:TYR:HD1	1:A:225:TRP:HB3	1.86	0.41
1:B:141:LEU:HA	1:B:141:LEU:HD23	1.79	0.41
1:B:313:ASN:HB2	1:B:314:ARG:HH21	1.86	0.41
1:A:93:ARG:HH21	1:A:128:ASN:ND2	2.19	0.41
1:B:235:ASP:HB3	1:B:270:GLN:O	2.21	0.41
1:B:282:LEU:HD22	1:B:303:TYR:OH	2.21	0.41
1:B:97:VAL:HG22	1:B:150:GLN:NE2	2.36	0.41
1:A:28:LYS:HB2	1:A:28:LYS:HE2	1.76	0.40
1:B:253:ASN:N	1:B:337:ASN:O	2.54	0.40
1:C:110:GLU:HB2	1:C:319:TYR:HE1	1.86	0.40
1:C:22:TYR:CD1	1:C:33:GLN:HG3	2.56	0.40
1:C:23:PHE:HB2	1:C:345:VAL:HB	2.02	0.40
1:A:334:TYR:O	1:A:338:GLY:N	2.53	0.40
1:C:86:TYR:CD1	1:C:86:TYR:N	2.89	0.40
1:A:155:ASN:OD1	1:A:169:ASN:ND2	2.42	0.40
1:C:125:GLY:HA3	2:C:420:HOH:O	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	351/356 (99%)	335 (95%)	16 (5%)	0	100	100
1	B	351/356 (99%)	337 (96%)	14 (4%)	0	100	100
1	C	351/356 (99%)	338 (96%)	13 (4%)	0	100	100
All	All	1053/1068 (99%)	1010 (96%)	43 (4%)	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	287/288 (100%)	279 (97%)	8 (3%)	43	69
1	B	287/288 (100%)	282 (98%)	5 (2%)	60	81
1	C	287/288 (100%)	278 (97%)	9 (3%)	40	66
All	All	861/864 (100%)	839 (97%)	22 (3%)	46	72

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	135	ASN
1	A	173	ASP
1	A	194	MET
1	A	204	SER
1	A	291	ASN
1	A	321	ASP
1	A	322	TYR
1	A	340	SER
1	B	173	ASP
1	B	174	SER
1	B	187	SER
1	B	284	SER
1	B	314	ARG
1	C	36	LEU
1	C	90	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	168	ASN
1	C	190	TYR
1	C	194	MET
1	C	243	SER
1	C	248	MET
1	C	276	ARG
1	C	322	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	261	GLN
1	A	352	GLN
1	B	150	GLN
1	B	261	GLN
1	B	268	GLN
1	B	352	GLN
1	C	150	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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, 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	353/356 (99%)	0.38	16 (4%) 33 26	32, 50, 80, 92	0
1	B	353/356 (99%)	0.67	39 (11%) 5 3	41, 64, 87, 108	0
1	C	353/356 (99%)	0.31	12 (3%) 45 38	32, 48, 70, 88	0
All	All	1059/1068 (99%)	0.46	67 (6%) 20 15	32, 53, 82, 108	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	165	VAL	7.6
1	A	294	ARG	5.8
1	B	86	TYR	5.7
1	B	164	ASN	5.0
1	B	286	GLY	4.7
1	A	86	TYR	4.5
1	C	291	ASN	4.4
1	B	289	LEU	4.0
1	B	291	ASN	4.0
1	B	160	ALA	3.9
1	B	314	ARG	3.8
1	B	235	ASP	3.6
1	A	219	GLY	3.4
1	C	158	GLN	3.4
1	B	330	ASN	3.3
1	B	296	ALA	3.3
1	B	72	ALA	3.3
1	A	138	PHE	3.0
1	B	234	ASN	3.0
1	A	218	ARG	3.0
1	B	233	ALA	3.0
1	C	141	LEU	2.9
1	C	219	GLY	2.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	220	ASP	2.9
1	B	273	PHE	2.8
1	C	60	VAL	2.8
1	B	269	TYR	2.8
1	B	285	LYS	2.7
1	B	162	ASP	2.6
1	B	190	TYR	2.6
1	B	161	GLN	2.6
1	B	272	ASP	2.6
1	B	184	PHE	2.5
1	B	297	ASP	2.5
1	C	83	VAL	2.4
1	B	295	TYR	2.4
1	A	339	ILE	2.4
1	A	176	VAL	2.4
1	C	184	PHE	2.4
1	B	298	LYS	2.4
1	B	342	ASP	2.3
1	B	344	ILE	2.3
1	B	58	TYR	2.3
1	A	72	ALA	2.3
1	B	141	LEU	2.3
1	A	118	TYR	2.3
1	B	261	GLN	2.3
1	B	340	SER	2.3
1	B	87	GLY	2.3
1	B	275	LEU	2.3
1	C	160	ALA	2.2
1	C	35	TYR	2.2
1	C	61	GLN	2.2
1	B	112	GLY	2.1
1	A	290	TYR	2.1
1	B	74	THR	2.1
1	A	127	ALA	2.1
1	A	287	LYS	2.1
1	C	81	ILE	2.1
1	A	129	GLY	2.1
1	B	163	ILE	2.1
1	A	259	LYS	2.1
1	B	59	ASN	2.1
1	A	158	GLN	2.0
1	A	4	TYR	2.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	300	LEU	2.0
1	B	142	VAL	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.