Full wwPDB X-ray Structure Validation Report (i)

Oct 9, 2023 – 08:22 AM EDT

PDB ID	:	7JQZ
Title	:	Crystal structure of Cfl2 wild-type from Burkholderia cenocepacia
Authors	:	Taher, N.M.; Madden, D.R.
Deposited on	:	2020-08-11
Resolution	:	2.20 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 (i) 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 (1)) were used in the production of this report:

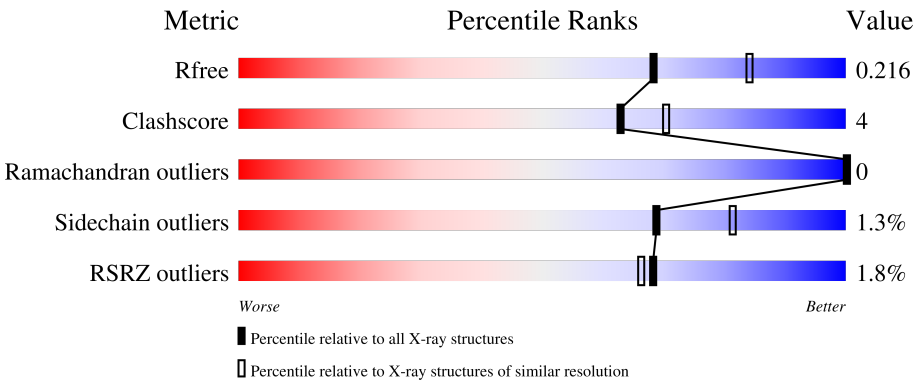
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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 <=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	309	.% 89% 6% • 5%
1	B	309	2% 85% 9% 5%
1	C	309	.% 83% 12% • 5%
1	D	309	.% 83% 11% 5%
1	E	309	4% 82% 13% • 5%

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Mol	Chain	Length	Quality of chain
1	F	309	84% 10% 5%
1	G	309	2% 86% 8% 5%
1	H	309	% 84% 11% 5%
1	I	309	3% 85% 10% 5%
1	J	309	2% 91% 5% 5%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 24478 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 Alpha/beta hydrolase fold.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	294	2348	1505	415	424	4	0	2	0
1	B	294	2337	1499	411	423	4	0	1	0
1	C	294	2326	1490	410	422	4	0	0	0
1	D	294	2326	1490	410	422	4	0	0	0
1	E	294	2326	1490	410	422	4	0	0	0
1	F	294	2337	1499	411	423	4	0	1	0
1	G	294	2337	1499	411	423	4	0	1	0
1	H	294	2337	1499	411	423	4	0	1	0
1	I	294	2326	1490	410	422	4	0	0	0
1	J	294	2337	1499	411	423	4	0	1	0

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	119	119	119	0	0
2	B	143	143	143	0	0
2	C	122	122	122	0	0
2	D	102	102	102	0	0

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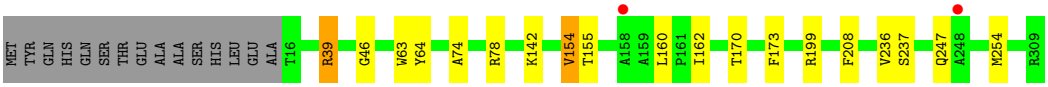
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	E	96	96	96	0	0
			Total	O		
2	F	137	137	137	0	0
			Total	O		
2	G	98	98	98	0	0
			Total	O		
2	H	119	119	119	0	0
			Total	O		
2	I	87	87	87	0	0
			Total	O		
2	J	118	118	118	0	0

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.

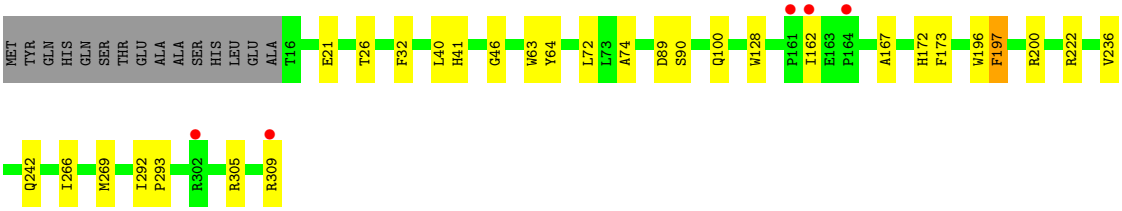
• Molecule 1: Alpha/beta hydrolase fold

Chain A:



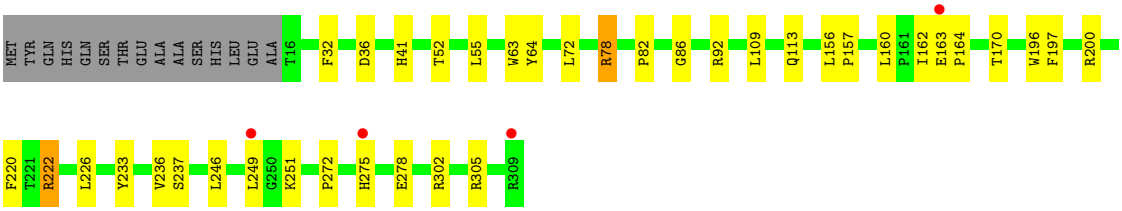
• Molecule 1: Alpha/beta hydrolase fold

Chain B:



• Molecule 1: Alpha/beta hydrolase fold

Chain C:

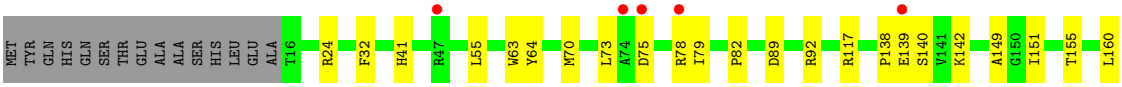


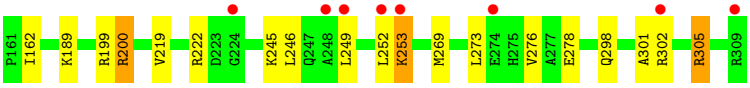
• Molecule 1: Alpha/beta hydrolase fold

Chain D:

• Molecule 1: Alpha/beta hydrolase fold

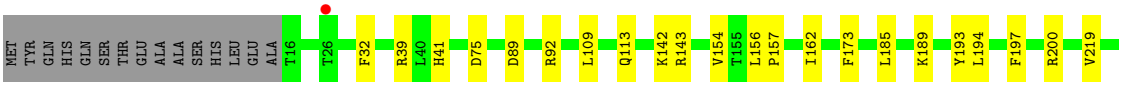
Chain E:





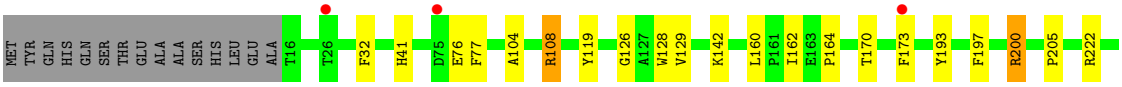
• Molecule 1: Alpha/beta hydrolase fold

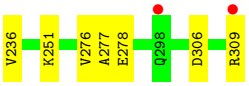
Chain F:



• Molecule 1: Alpha/beta hydrolase fold

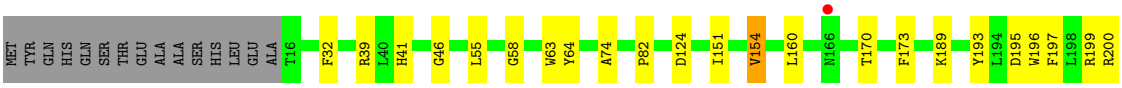
Chain G:

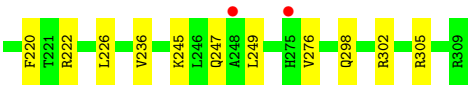




• Molecule 1: Alpha/beta hydrolase fold

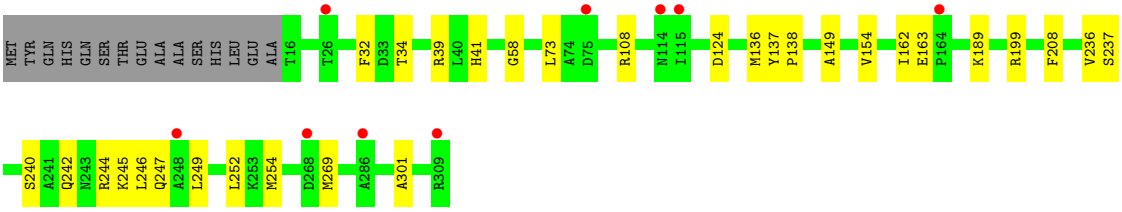
Chain H:





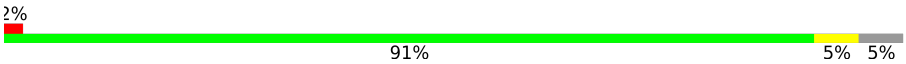
• Molecule 1: Alpha/beta hydrolase fold

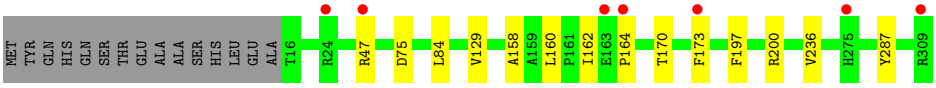
Chain I:



- Molecule 1: Alpha/beta hydrolase fold

Chain J:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	182.73Å 210.47Å 86.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.54 – 2.20 48.54 – 2.20	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.54-2.20) 100.0 (48.54-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.11 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.185 , 0.216 0.186 , 0.216	Depositor DCC
R_{free} test set	8500 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å^2)	37.3	Xtriage
Anisotropy	0.252	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 37.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	24478	wwPDB-VP
Average B, all atoms (Å^2)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.01% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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	0.54	0/2414	0.66	0/3285
1	B	0.51	0/2403	0.64	0/3271
1	C	0.52	0/2391	0.68	0/3255
1	D	0.54	0/2391	0.70	0/3255
1	E	0.54	0/2391	0.68	0/3255
1	F	0.50	0/2403	0.64	0/3271
1	G	0.53	0/2403	0.62	0/3271
1	H	0.55	0/2403	0.66	0/3271
1	I	0.52	0/2391	0.67	0/3255
1	J	0.53	0/2403	0.65	0/3271
All	All	0.53	0/23993	0.66	0/32660

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	2348	0	2287	16	0
1	B	2337	0	2275	19	0
1	C	2326	0	2267	27	0
1	D	2326	0	2267	24	0
1	E	2326	0	2267	33	0
1	F	2337	0	2275	27	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	G	2337	0	2275	18	0
1	H	2337	0	2275	21	0
1	I	2326	0	2267	20	0
1	J	2337	0	2275	9	0
2	A	119	0	0	1	0
2	B	143	0	0	0	0
2	C	122	0	0	0	0
2	D	102	0	0	2	0
2	E	96	0	0	0	0
2	F	137	0	0	1	0
2	G	98	0	0	0	0
2	H	119	0	0	2	0
2	I	87	0	0	0	0
2	J	118	0	0	1	0
All	All	24478	0	22730	197	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 4.

All (197) 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:I:149:ALA:H	1:I:269:MET:HE1	1.57	0.69
1:H:247:GLN:HG3	1:H:276:VAL:HG12	1.74	0.69
1:D:160:LEU:HB3	1:D:161:PRO:HD2	1.75	0.68
1:F:154:VAL:HG13	1:F:276:VAL:HG11	1.76	0.68
1:F:189:LYS:NZ	2:F:402:HOH:O	2.26	0.68
1:I:237:SER:HB3	1:J:164:PRO:HD2	1.74	0.67
1:D:160:LEU:HD13	1:D:170:THR:HB	1.77	0.67
1:E:138:PRO:HB2	1:E:253:LYS:HE3	1.77	0.66
1:F:154:VAL:CG1	1:F:276:VAL:HG11	2.26	0.66
1:I:136:MET:C	1:I:138:PRO:HD3	2.15	0.66
1:C:237:SER:HB3	1:D:164:PRO:HD2	1.77	0.65
1:C:272:PRO:O	1:C:275:HIS:CE1	2.50	0.65
1:E:149:ALA:H	1:E:269:MET:HE1	1.61	0.65
1:A:173[B]:PHE:HZ	1:A:236:VAL:HG21	1.63	0.64
1:H:302:ARG:HH12	1:H:305:ARG:HH11	1.45	0.64
1:C:272:PRO:O	1:C:275:HIS:ND1	2.31	0.63
1:A:154:VAL:HG23	1:A:247:GLN:OE1	1.98	0.63
1:B:173[B]:PHE:HZ	1:B:236:VAL:HG21	1.62	0.63
1:D:237:SER:HB2	2:D:426:HOH:O	1.97	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:163:GLU:OE1	1:C:163:GLU:N	2.24	0.62
1:I:199:ARG:NH1	1:I:208:PHE:O	2.34	0.60
1:F:298:GLN:O	1:F:302:ARG:HG2	2.01	0.60
1:G:251:LYS:HB2	1:G:276:VAL:C	2.22	0.60
1:D:247:GLN:O	1:D:247:GLN:HG3	2.01	0.59
1:I:137:TYR:N	1:I:138:PRO:HD3	2.16	0.58
1:F:75:ASP:OD1	1:F:305:ARG:NH2	2.35	0.58
1:I:244:ARG:O	1:I:247:GLN:HB3	2.03	0.58
1:A:78:ARG:HH11	1:A:78:ARG:HG2	1.68	0.58
1:H:173[B]:PHE:HZ	1:H:236:VAL:HG21	1.67	0.58
1:J:173[B]:PHE:HZ	1:J:236:VAL:HG21	1.69	0.58
1:C:163:GLU:OE2	1:D:158:ALA:O	2.22	0.58
1:E:245:LYS:O	1:E:249:LEU:HD13	2.04	0.58
1:G:76:GLU:OE2	1:G:309:ARG:NH1	2.37	0.58
1:H:151:ILE:HB	1:H:154:VAL:HG23	1.85	0.57
1:D:199:ARG:NH1	1:D:208:PHE:O	2.36	0.57
1:A:154:VAL:HG13	1:A:155:THR:HG23	1.86	0.56
1:E:92:ARG:HH21	1:E:222:ARG:HG3	1.69	0.56
1:E:298:GLN:O	1:E:302:ARG:HG3	2.06	0.56
1:E:149:ALA:N	1:E:269:MET:HE1	2.21	0.55
1:C:160:LEU:HD13	1:C:170:THR:HB	1.89	0.55
1:C:92:ARG:NH2	1:C:222:ARG:CD	2.70	0.55
1:E:89:ASP:HA	1:E:222:ARG:HH11	1.72	0.54
1:E:92:ARG:NH2	1:E:222:ARG:CG	2.71	0.54
1:I:73:LEU:HD21	1:I:301:ALA:HB1	1.89	0.54
1:I:236:VAL:O	1:I:240:SER:HB2	2.06	0.54
1:B:305:ARG:HG2	1:B:309:ARG:HH22	1.72	0.54
1:A:173[B]:PHE:CZ	1:A:236:VAL:HG21	2.43	0.54
1:C:92:ARG:NH2	1:C:222:ARG:HD2	2.23	0.54
1:A:162:ILE:HB	1:B:173[B]:PHE:CD2	2.43	0.53
1:E:117:ARG:CZ	1:E:142:LYS:HG2	2.39	0.53
1:E:199:ARG:O	1:H:39:ARG:NH2	2.42	0.53
1:F:143:ARG:HD3	1:F:255:PRO:HG2	1.90	0.52
1:H:160:LEU:HD13	1:H:170:THR:HB	1.91	0.52
1:B:89:ASP:HA	1:B:222:ARG:HH11	1.73	0.52
1:E:139:GLU:HG2	1:E:140:SER:N	2.25	0.52
1:I:254:MET:SD	1:I:254:MET:N	2.82	0.52
1:C:163:GLU:OE2	1:D:158:ALA:HB1	2.09	0.52
1:E:32:PHE:CE2	1:E:41:HIS:HB2	2.45	0.52
1:A:199:ARG:NH1	1:A:208:PHE:O	2.43	0.51
1:E:92:ARG:HH21	1:E:222:ARG:CG	2.24	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:55:LEU:O	1:E:82:PRO:HD2	2.11	0.51
1:B:89:ASP:HA	1:B:222:ARG:NH1	2.26	0.51
1:H:189:LYS:NZ	2:H:401:HOH:O	2.35	0.50
1:F:173[C]:PHE:HZ	1:F:236:VAL:HG21	1.76	0.50
1:F:257:LEU:HB2	1:F:307:PHE:CE1	2.46	0.50
1:A:39:ARG:HH22	1:D:200:ARG:HA	1.77	0.50
1:F:302:ARG:HH22	1:F:305:ARG:NH1	2.09	0.50
1:C:92:ARG:HH21	1:C:222:ARG:HD3	1.76	0.50
1:C:233:TYR:HA	1:C:236:VAL:HG13	1.92	0.50
1:B:100:GLN:HB2	1:B:242:GLN:OE1	2.13	0.49
1:E:138:PRO:CB	1:E:253:LYS:HE3	2.40	0.49
1:C:251:LYS:HE2	1:C:278:GLU:HA	1.94	0.49
1:C:220:PHE:CZ	1:C:226:LEU:HD13	2.48	0.49
1:G:193:TYR:CZ	1:G:197:PHE:HE1	2.30	0.48
1:F:302:ARG:NH2	1:F:305:ARG:HH11	2.12	0.48
1:D:154:VAL:HG12	1:D:155:THR:HG23	1.95	0.48
1:E:73:LEU:HB2	1:E:79:ILE:HD11	1.95	0.48
1:F:32:PHE:CE2	1:F:41:HIS:HB2	2.47	0.48
1:D:163:GLU:HG3	1:D:166:ASN:HB2	1.96	0.48
1:E:273:LEU:O	1:E:276:VAL:HG22	2.14	0.47
1:D:222:ARG:O	1:D:222:ARG:HG3	2.10	0.47
1:G:160:LEU:HD13	1:G:170:THR:HB	1.95	0.47
1:G:251:LYS:HB2	1:G:276:VAL:O	2.14	0.47
1:H:298:GLN:O	1:H:302:ARG:HG2	2.14	0.47
1:I:154:VAL:HG12	1:I:247:GLN:NE2	2.29	0.47
1:B:162:ILE:O	1:B:167:ALA:HB2	2.14	0.47
1:J:160:LEU:HD13	1:J:170:THR:HB	1.96	0.47
1:A:78:ARG:HG2	1:A:78:ARG:NH1	2.29	0.47
1:F:156:LEU:HD21	1:F:173[C]:PHE:CE1	2.49	0.47
1:A:237:SER:HB2	2:A:468:HOH:O	2.15	0.47
1:F:302:ARG:HH22	1:F:305:ARG:HH11	1.62	0.47
1:C:32:PHE:CE2	1:C:41:HIS:HB2	2.49	0.47
1:H:195:ASP:OD1	1:H:199:ARG:NH1	2.45	0.47
1:E:89:ASP:HA	1:E:222:ARG:NH1	2.29	0.47
1:E:189:LYS:HD3	1:F:185:LEU:HA	1.96	0.47
1:I:34:THR:OG1	1:I:39:ARG:HG2	2.15	0.47
1:B:40:LEU:HD23	1:B:90:SER:HB3	1.97	0.46
1:C:52:THR:HA	1:C:78:ARG:O	2.15	0.46
1:D:100:GLN:HB2	1:D:242:GLN:OE1	2.15	0.46
1:G:173[A]:PHE:HZ	1:G:236:VAL:HG21	1.80	0.46
1:H:193:TYR:CZ	1:H:197:PHE:HE1	2.34	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:196:TRP:O	1:C:200:ARG:HB3	2.16	0.46
1:C:55:LEU:O	1:C:82:PRO:HD2	2.16	0.46
1:C:163:GLU:H	1:C:163:GLU:CD	2.17	0.46
1:G:162:ILE:HG21	1:H:160:LEU:HD12	1.97	0.45
1:I:32:PHE:CE2	1:I:41:HIS:HB2	2.51	0.45
1:C:164:PRO:HD2	1:D:237:SER:OG	2.15	0.45
1:C:246:LEU:O	1:C:249:LEU:HB2	2.17	0.45
1:D:113:GLN:HB2	1:D:115:ILE:HD12	1.97	0.45
1:E:75:ASP:OD1	1:E:305:ARG:NH2	2.49	0.45
1:E:151:ILE:H	1:E:155:THR:HB	1.80	0.45
1:G:251:LYS:HD3	1:G:277:ALA:C	2.37	0.45
1:B:266:ILE:HG21	1:B:269:MET:HG3	1.99	0.44
1:B:46:GLY:HA2	1:B:74:ALA:HB1	2.00	0.44
1:B:63:TRP:CG	1:B:64:TYR:N	2.85	0.44
1:C:72:LEU:O	1:C:305:ARG:NH2	2.49	0.44
1:B:72:LEU:O	1:B:305:ARG:NH2	2.50	0.44
1:B:196:TRP:O	1:B:200:ARG:HB3	2.17	0.44
1:E:246:LEU:O	1:E:249:LEU:HB2	2.17	0.44
1:D:47:ARG:HD2	1:D:75:ASP:O	2.17	0.44
1:C:92:ARG:NH2	1:C:222:ARG:HD3	2.32	0.44
1:E:70:MET:HG2	1:E:79:ILE:HG21	1.99	0.44
1:F:193:TYR:CZ	1:F:197:PHE:HE1	2.36	0.44
1:D:32:PHE:CE2	1:D:41:HIS:HB2	2.53	0.44
1:J:84:LEU:HD13	1:J:129:VAL:HG21	1.99	0.44
1:H:245:LYS:O	1:H:249:LEU:HG	2.18	0.44
1:J:173[B]:PHE:CZ	1:J:236:VAL:HG21	2.52	0.44
1:A:142:LYS:HE2	1:A:142:LYS:HB3	1.89	0.43
1:D:73:LEU:HB2	1:D:79:ILE:HD11	2.00	0.43
1:F:265:SER:HB2	1:F:266:ILE:HD12	2.00	0.43
1:C:109:LEU:O	1:C:113:GLN:HG2	2.17	0.43
1:E:117:ARG:NH2	1:E:142:LYS:HG2	2.33	0.43
1:E:219:VAL:O	1:E:222:ARG:HG2	2.18	0.43
1:I:242:GLN:O	1:I:246:LEU:HG	2.18	0.43
1:G:76:GLU:HG2	1:G:77:PHE:CD2	2.53	0.43
1:B:21:GLU:OE1	1:B:26:THR:HA	2.18	0.43
1:C:156:LEU:HD12	1:C:157:PRO:HD2	2.01	0.43
1:H:220:PHE:CZ	1:H:226:LEU:HD13	2.53	0.43
1:I:162:ILE:HG12	1:J:158:ALA:O	2.18	0.43
1:B:128:TRP:HZ2	1:B:173[A]:PHE:HE1	1.67	0.43
1:H:58:GLY:HA3	1:H:124:ASP:HB3	2.00	0.43
1:A:63:TRP:CG	1:A:64:TYR:N	2.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:160:LEU:HD13	1:A:170:THR:HB	2.01	0.42
1:I:58:GLY:HA3	1:I:124:ASP:HB3	2.01	0.42
1:A:142:LYS:O	1:A:254:MET:HB2	2.19	0.42
1:D:63:TRP:HA	2:D:460:HOH:O	2.20	0.42
1:I:163:GLU:H	1:I:163:GLU:HG2	1.64	0.42
1:J:287:TYR:HA	2:J:476:HOH:O	2.19	0.42
1:D:55:LEU:HD23	1:D:121:ALA:HB3	2.02	0.42
1:F:156:LEU:HA	1:F:157:PRO:HD3	1.92	0.42
1:F:89:ASP:HA	1:F:222:ARG:HH11	1.85	0.42
1:A:46:GLY:HA2	1:A:74:ALA:HB1	2.02	0.42
1:E:92:ARG:NH2	1:E:222:ARG:HG3	2.33	0.42
1:B:32:PHE:CE2	1:B:41:HIS:HB2	2.53	0.42
1:D:55:LEU:O	1:D:82:PRO:HD2	2.18	0.42
1:G:104:ALA:O	1:G:108:ARG:HD2	2.20	0.42
1:G:306:ASP:O	1:G:309:ARG:HG2	2.19	0.42
1:E:200:ARG:HA	1:H:39:ARG:HH22	1.84	0.42
1:F:39:ARG:HD2	1:G:205:PRO:HD3	2.00	0.42
1:G:251:LYS:HG3	1:G:278:GLU:HG2	2.01	0.42
1:J:47:ARG:HD2	1:J:75:ASP:O	2.20	0.42
1:I:252:LEU:HD23	1:I:252:LEU:HA	1.84	0.42
1:G:126:GLY:O	1:G:129:VAL:HG22	2.19	0.41
1:B:172:HIS:CG	1:B:197:PHE:HZ	2.38	0.41
1:G:32:PHE:CE2	1:G:41:HIS:HB2	2.56	0.41
1:I:245:LYS:O	1:I:249:LEU:HD13	2.19	0.41
1:D:76:GLU:OE1	1:D:309:ARG:NH1	2.46	0.41
1:E:24:ARG:NH1	1:E:24:ARG:HG3	2.35	0.41
1:F:257:LEU:HB2	1:F:307:PHE:CD1	2.55	0.41
1:H:32:PHE:CE2	1:H:41:HIS:HB2	2.56	0.41
1:H:46:GLY:HA2	1:H:74:ALA:HB1	2.02	0.41
1:E:63:TRP:CG	1:E:64:TYR:N	2.87	0.41
1:H:196:TRP:O	1:H:200:ARG:HB3	2.20	0.41
1:B:292:ILE:HB	1:B:293:PRO:HD3	2.02	0.41
1:F:39:ARG:HH12	1:G:200:ARG:HA	1.85	0.41
1:G:128:TRP:CZ2	1:G:173[B]:PHE:HE1	2.39	0.41
1:H:220:PHE:HB2	2:H:442:HOH:O	2.20	0.41
1:A:160:LEU:HD12	1:B:162:ILE:HG21	2.02	0.41
1:D:75:ASP:OD1	1:D:76:GLU:N	2.54	0.41
1:E:162:ILE:HD12	1:E:162:ILE:C	2.41	0.41
1:F:142:LYS:O	1:F:143:ARG:HG2	2.21	0.41
1:I:137:TYR:N	1:I:138:PRO:CD	2.81	0.41
1:I:244:ARG:HG2	1:I:247:GLN:OE1	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:301:ALA:O	1:E:305:ARG:HB2	2.20	0.41
1:F:92:ARG:NH2	1:F:222:ARG:HD3	2.36	0.41
1:F:194:LEU:HD23	1:F:194:LEU:HA	1.83	0.41
1:H:55:LEU:O	1:H:82:PRO:HD2	2.21	0.41
1:E:160:LEU:HG	1:F:162:ILE:HG21	2.04	0.40
1:F:109:LEU:O	1:F:113:GLN:HG2	2.21	0.40
1:E:252:LEU:O	1:E:278:GLU:HB2	2.21	0.40
1:J:162:ILE:O	1:J:162:ILE:HG13	2.20	0.40
1:C:63:TRP:CG	1:C:64:TYR:N	2.89	0.40
1:C:162:ILE:HG21	1:D:160:LEU:CD1	2.51	0.40
1:F:219:VAL:O	1:F:222:ARG:HG2	2.22	0.40
1:G:119:TYR:OH	1:G:142:LYS:HE3	2.22	0.40
1:C:86:GLY:HA2	1:C:92:ARG:HG3	2.04	0.40
1:H:63:TRP:CG	1:H:64:TYR:N	2.88	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	294/309 (95%)	287 (98%)	7 (2%)	0	100	100
1	B	293/309 (95%)	285 (97%)	8 (3%)	0	100	100
1	C	292/309 (94%)	286 (98%)	6 (2%)	0	100	100
1	D	292/309 (94%)	286 (98%)	6 (2%)	0	100	100
1	E	292/309 (94%)	284 (97%)	8 (3%)	0	100	100
1	F	293/309 (95%)	285 (97%)	8 (3%)	0	100	100
1	G	293/309 (95%)	284 (97%)	9 (3%)	0	100	100
1	H	293/309 (95%)	286 (98%)	7 (2%)	0	100	100
1	I	292/309 (94%)	283 (97%)	9 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	J	293/309 (95%)	286 (98%)	7 (2%)	0	100	100
All	All	2927/3090 (95%)	2852 (97%)	75 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	235/245 (96%)	233 (99%)	2 (1%)	78	88
1	B	234/245 (96%)	233 (100%)	1 (0%)	91	96
1	C	233/245 (95%)	228 (98%)	5 (2%)	53	67
1	D	233/245 (95%)	227 (97%)	6 (3%)	46	58
1	E	233/245 (95%)	229 (98%)	4 (2%)	60	74
1	F	234/245 (96%)	232 (99%)	2 (1%)	78	88
1	G	234/245 (96%)	230 (98%)	4 (2%)	60	74
1	H	234/245 (96%)	232 (99%)	2 (1%)	78	88
1	I	233/245 (95%)	231 (99%)	2 (1%)	78	88
1	J	234/245 (96%)	232 (99%)	2 (1%)	78	88
All	All	2337/2450 (95%)	2307 (99%)	30 (1%)	69	81

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	ARG
1	A	154	VAL
1	B	197	PHE
1	C	36	ASP
1	C	78	ARG
1	C	197	PHE
1	C	222	ARG
1	C	302	ARG

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Mol	Chain	Res	Type
1	D	36	ASP
1	D	85	PRO
1	D	139	GLU
1	D	223	ASP
1	D	237	SER
1	D	253	LYS
1	E	78	ARG
1	E	200	ARG
1	E	253	LYS
1	E	305	ARG
1	F	200	ARG
1	F	302	ARG
1	G	108	ARG
1	G	164	PRO
1	G	200	ARG
1	G	222	ARG
1	H	154	VAL
1	H	222	ARG
1	I	108	ARG
1	I	189	LYS
1	J	197	PHE
1	J	200	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 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 (i)

6.1 Protein, DNA and RNA chains (i)

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	294/309 (95%)	-0.25	2 (0%)	87	86	24, 34, 51, 61	0
1	B	294/309 (95%)	-0.25	5 (1%)	70	68	25, 34, 51, 74	0
1	C	294/309 (95%)	-0.17	4 (1%)	75	73	24, 37, 57, 73	0
1	D	294/309 (95%)	-0.01	3 (1%)	82	81	26, 39, 59, 73	0
1	E	294/309 (95%)	0.07	13 (4%)	34	32	25, 44, 62, 74	0
1	F	294/309 (95%)	-0.31	1 (0%)	94	93	25, 35, 52, 64	0
1	G	294/309 (95%)	-0.13	5 (1%)	70	68	24, 37, 56, 73	0
1	H	294/309 (95%)	-0.26	3 (1%)	82	81	23, 34, 55, 68	0
1	I	294/309 (95%)	0.05	9 (3%)	49	47	26, 44, 70, 83	0
1	J	294/309 (95%)	-0.17	7 (2%)	59	56	25, 35, 55, 82	0
All	All	2940/3090 (95%)	-0.14	52 (1%)	68	66	23, 37, 59, 83	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	26	THR	3.7
1	J	164	PRO	3.5
1	H	248	ALA	3.3
1	E	75	ASP	3.1
1	I	114	ASN	3.1
1	E	47	ARG	3.0
1	H	166	ASN	2.9
1	E	248	ALA	2.9
1	G	309	ARG	2.9
1	C	309	ARG	2.8
1	A	248	ALA	2.8
1	I	248	ALA	2.8
1	I	286	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	302	ARG	2.7
1	C	163	GLU	2.7
1	G	173[A]	PHE	2.7
1	I	268	ASP	2.7
1	E	249	LEU	2.6
1	E	253	LYS	2.6
1	E	309	ARG	2.6
1	B	302	ARG	2.6
1	I	75	ASP	2.6
1	I	309	ARG	2.5
1	I	115	ILE	2.5
1	J	309	ARG	2.4
1	J	275	HIS	2.4
1	B	164	PRO	2.4
1	B	309	ARG	2.4
1	B	162	ILE	2.4
1	C	249	LEU	2.4
1	I	26	THR	2.4
1	G	298	GLN	2.4
1	J	173[A]	PHE	2.4
1	F	26	THR	2.3
1	G	75	ASP	2.3
1	A	158	ALA	2.3
1	D	298	GLN	2.3
1	H	275	HIS	2.3
1	E	139	GLU	2.2
1	E	224	GLY	2.2
1	J	24	ARG	2.2
1	D	164	PRO	2.2
1	I	164	PRO	2.2
1	J	47	ARG	2.2
1	B	161	PRO	2.2
1	E	274	GLU	2.1
1	E	252	LEU	2.1
1	C	275	HIS	2.1
1	E	74	ALA	2.1
1	E	78	ARG	2.0
1	E	302	ARG	2.0
1	J	163	GLU	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.