



Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 01:52 pm BST

PDB ID : 2J7V
Title : Structure of PBP-A
Authors : Evrard, C.; Declercq, J.P.
Deposited on : 2006-10-17
Resolution : 1.90 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

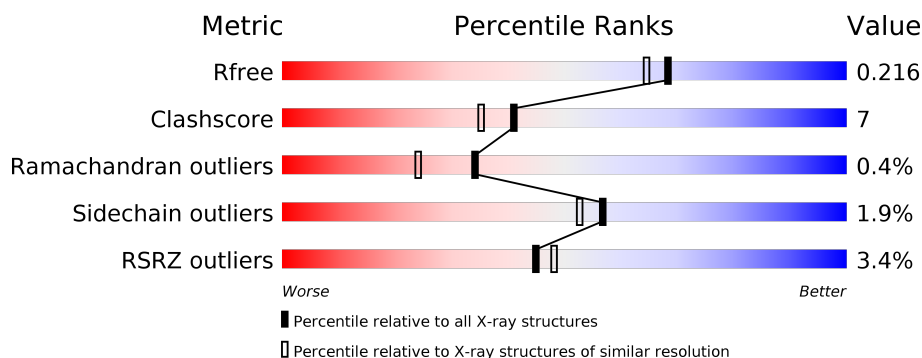
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria 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	298	<div> <div>2%</div> <div> <div></div> <div>80%</div> <div>8%</div> <div>11%</div> </div> </div>
1	B	298	<div> <div>%</div> <div> <div></div> <div>80%</div> <div>9%</div> <div>11%</div> </div> </div>
1	C	298	<div> <div>4%</div> <div> <div></div> <div>77%</div> <div>11%</div> <div>9%</div> </div> </div>
1	D	298	<div> <div>5%</div> <div> <div></div> <div>77%</div> <div>14%</div> <div>8%</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9125 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 TLL2115 PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	265	Total	C	N	O	S	0	0	0
			2041	1282	361	387	11			
1	B	265	Total	C	N	O	S	0	0	0
			2040	1281	361	387	11			
1	C	270	Total	C	N	O	S	0	0	0
			2072	1300	366	395	11			
1	D	274	Total	C	N	O	S	0	0	0
			2102	1320	370	400	12			

There are 88 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP Q8DH45
A	278	GLU	-	expression tag	UNP Q8DH45
A	279	GLN	-	expression tag	UNP Q8DH45
A	280	LYS	-	expression tag	UNP Q8DH45
A	281	LEU	-	expression tag	UNP Q8DH45
A	282	ILE	-	expression tag	UNP Q8DH45
A	283	SER	-	expression tag	UNP Q8DH45
A	284	GLU	-	expression tag	UNP Q8DH45
A	285	GLU	-	expression tag	UNP Q8DH45
A	286	ASP	-	expression tag	UNP Q8DH45
A	287	LEU	-	expression tag	UNP Q8DH45
A	288	ASN	-	expression tag	UNP Q8DH45
A	289	SER	-	expression tag	UNP Q8DH45
A	290	ALA	-	expression tag	UNP Q8DH45
A	291	VAL	-	expression tag	UNP Q8DH45
A	292	ASP	-	expression tag	UNP Q8DH45
A	293	HIS	-	expression tag	UNP Q8DH45
A	294	HIS	-	expression tag	UNP Q8DH45
A	295	HIS	-	expression tag	UNP Q8DH45
A	296	HIS	-	expression tag	UNP Q8DH45
A	297	HIS	-	expression tag	UNP Q8DH45

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Chain	Residue	Modelled	Actual	Comment	Reference
A	298	HIS	-	expression tag	UNP Q8DH45
B	1	MET	-	expression tag	UNP Q8DH45
B	278	GLU	-	expression tag	UNP Q8DH45
B	279	GLN	-	expression tag	UNP Q8DH45
B	280	LYS	-	expression tag	UNP Q8DH45
B	281	LEU	-	expression tag	UNP Q8DH45
B	282	ILE	-	expression tag	UNP Q8DH45
B	283	SER	-	expression tag	UNP Q8DH45
B	284	GLU	-	expression tag	UNP Q8DH45
B	285	GLU	-	expression tag	UNP Q8DH45
B	286	ASP	-	expression tag	UNP Q8DH45
B	287	LEU	-	expression tag	UNP Q8DH45
B	288	ASN	-	expression tag	UNP Q8DH45
B	289	SER	-	expression tag	UNP Q8DH45
B	290	ALA	-	expression tag	UNP Q8DH45
B	291	VAL	-	expression tag	UNP Q8DH45
B	292	ASP	-	expression tag	UNP Q8DH45
B	293	HIS	-	expression tag	UNP Q8DH45
B	294	HIS	-	expression tag	UNP Q8DH45
B	295	HIS	-	expression tag	UNP Q8DH45
B	296	HIS	-	expression tag	UNP Q8DH45
B	297	HIS	-	expression tag	UNP Q8DH45
B	298	HIS	-	expression tag	UNP Q8DH45
C	1	MET	-	expression tag	UNP Q8DH45
C	278	GLU	-	expression tag	UNP Q8DH45
C	279	GLN	-	expression tag	UNP Q8DH45
C	280	LYS	-	expression tag	UNP Q8DH45
C	281	LEU	-	expression tag	UNP Q8DH45
C	282	ILE	-	expression tag	UNP Q8DH45
C	283	SER	-	expression tag	UNP Q8DH45
C	284	GLU	-	expression tag	UNP Q8DH45
C	285	GLU	-	expression tag	UNP Q8DH45
C	286	ASP	-	expression tag	UNP Q8DH45
C	287	LEU	-	expression tag	UNP Q8DH45
C	288	ASN	-	expression tag	UNP Q8DH45
C	289	SER	-	expression tag	UNP Q8DH45
C	290	ALA	-	expression tag	UNP Q8DH45
C	291	VAL	-	expression tag	UNP Q8DH45
C	292	ASP	-	expression tag	UNP Q8DH45
C	293	HIS	-	expression tag	UNP Q8DH45
C	294	HIS	-	expression tag	UNP Q8DH45
C	295	HIS	-	expression tag	UNP Q8DH45

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Chain	Residue	Modelled	Actual	Comment	Reference
C	296	HIS	-	expression tag	UNP Q8DH45
C	297	HIS	-	expression tag	UNP Q8DH45
C	298	HIS	-	expression tag	UNP Q8DH45
D	1	MET	-	expression tag	UNP Q8DH45
D	278	GLU	-	expression tag	UNP Q8DH45
D	279	GLN	-	expression tag	UNP Q8DH45
D	280	LYS	-	expression tag	UNP Q8DH45
D	281	LEU	-	expression tag	UNP Q8DH45
D	282	ILE	-	expression tag	UNP Q8DH45
D	283	SER	-	expression tag	UNP Q8DH45
D	284	GLU	-	expression tag	UNP Q8DH45
D	285	GLU	-	expression tag	UNP Q8DH45
D	286	ASP	-	expression tag	UNP Q8DH45
D	287	LEU	-	expression tag	UNP Q8DH45
D	288	ASN	-	expression tag	UNP Q8DH45
D	289	SER	-	expression tag	UNP Q8DH45
D	290	ALA	-	expression tag	UNP Q8DH45
D	291	VAL	-	expression tag	UNP Q8DH45
D	292	ASP	-	expression tag	UNP Q8DH45
D	293	HIS	-	expression tag	UNP Q8DH45
D	294	HIS	-	expression tag	UNP Q8DH45
D	295	HIS	-	expression tag	UNP Q8DH45
D	296	HIS	-	expression tag	UNP Q8DH45
D	297	HIS	-	expression tag	UNP Q8DH45
D	298	HIS	-	expression tag	UNP Q8DH45

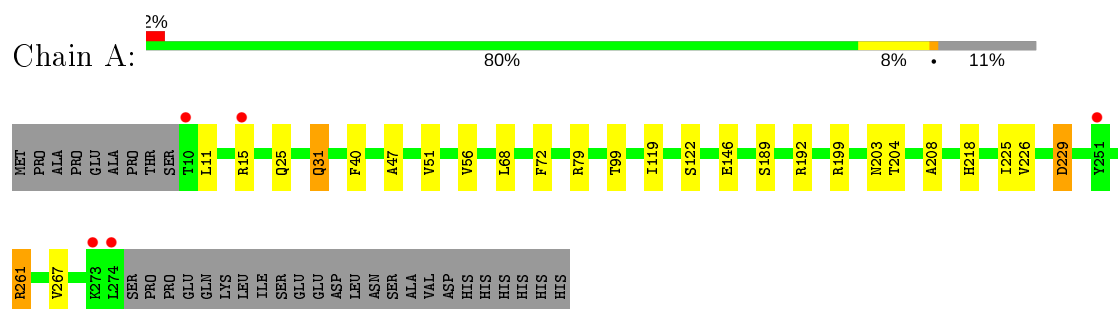
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	217	Total O 217 217	0	0
2	B	287	Total O 287 287	0	0
2	C	256	Total O 256 256	0	0
2	D	110	Total O 110 110	0	0

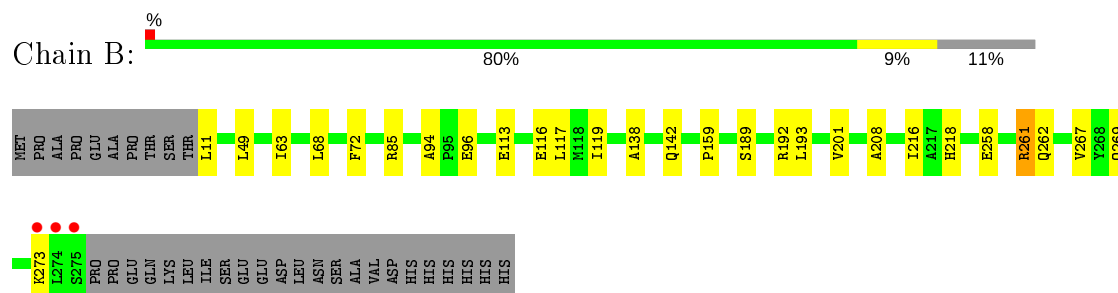
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

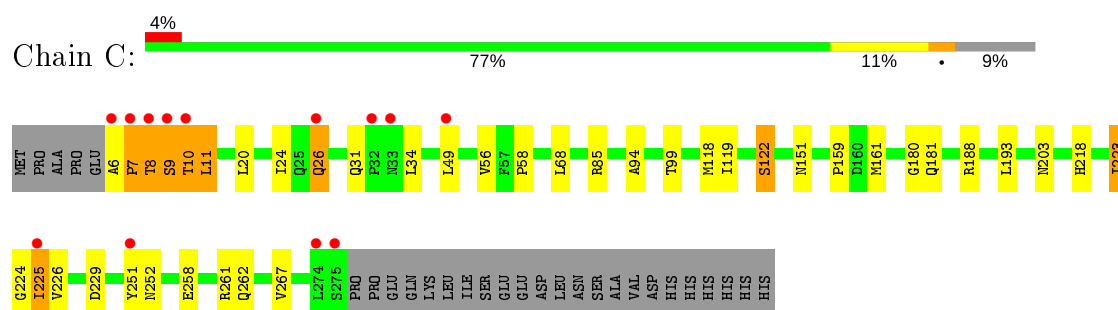
• Molecule 1: TLL2115 PROTEIN



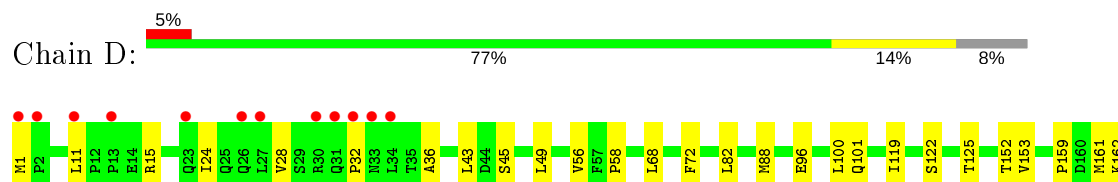
• Molecule 1: TLL2115 PROTEIN



• Molecule 1: TLL2115 PROTEIN



• Molecule 1: TLL2115 PROTEIN





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	87.79 Å 91.90 Å 147.27 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.20 – 1.90 39.24 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.8 (39.20-1.90) 99.8 (39.24-1.90)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.92 (at 1.89 Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.172 , 0.213 0.177 , 0.216	Depositor DCC
R_{free} test set	4726 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	21.7	Xtriage
Anisotropy	0.106	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 50.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.021 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9125	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.97% 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	0.70	0/2074	0.82	6/2814 (0.2%)
1	B	0.74	0/2073	0.80	2/2812 (0.1%)
1	C	0.74	0/2106	0.82	2/2859 (0.1%)
1	D	0.57	0/2138	0.69	1/2904 (0.0%)
All	All	0.69	0/8391	0.79	11/11389 (0.1%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	261	ARG	NE-CZ-NH2	-7.99	116.31	120.30
1	A	199	ARG	NE-CZ-NH1	7.85	124.23	120.30
1	C	261	ARG	NE-CZ-NH2	-7.56	116.52	120.30
1	A	199	ARG	NE-CZ-NH2	-7.49	116.56	120.30
1	A	229	ASP	CB-CG-OD1	6.58	124.22	118.30
1	D	188	ARG	NE-CZ-NH2	-5.98	117.31	120.30
1	A	261	ARG	NE-CZ-NH2	-5.93	117.33	120.30
1	A	229	ASP	CB-CG-OD2	-5.79	113.09	118.30
1	A	79	ARG	NE-CZ-NH2	-5.75	117.42	120.30
1	B	261	ARG	NE-CZ-NH1	5.48	123.04	120.30
1	C	261	ARG	NE-CZ-NH1	5.46	123.03	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2041	0	2078	20	0
1	B	2040	0	2076	19	1
1	C	2072	0	2107	42	0
1	D	2102	0	2139	39	1
2	A	217	0	0	4	0
2	B	287	0	0	4	0
2	C	256	0	0	7	0
2	D	110	0	0	4	0
All	All	9125	0	8400	118	1

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 7.

All (118) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:88:MET:CE	1:D:125:THR:HG21	1.58	1.32
1:D:88:MET:HE2	1:D:125:THR:HG21	1.22	1.14
1:D:88:MET:HE3	1:D:125:THR:HG21	1.37	1.03
1:C:20:LEU:CD2	1:C:49:LEU:HD11	1.94	0.97
1:C:20:LEU:HD23	1:C:49:LEU:HD11	1.51	0.92
1:D:88:MET:HE2	1:D:125:THR:CG2	2.04	0.87
1:D:166:THR:HG21	2:D:2069:HOH:O	1.79	0.82
1:B:49:LEU:CD1	1:B:267:VAL:HG22	2.10	0.80
1:D:153:VAL:HB	1:D:166:THR:HG22	1.63	0.80
1:D:96:GLU:HG3	1:D:161:MET:HE2	1.65	0.77
1:C:20:LEU:HD21	1:C:49:LEU:HD11	1.67	0.77
1:C:49:LEU:CD1	1:C:267:VAL:HG22	2.15	0.76
1:C:161:MET:HE1	1:C:223:ILE:HG22	1.69	0.74
1:D:43:LEU:HD22	1:D:177:LEU:HD13	1.71	0.72
1:C:68:LEU:HD22	1:C:119:ILE:HG13	1.70	0.71
1:C:6:ALA:HB3	1:C:7:PRO:HD3	1.73	0.71
1:B:85:ARG:HD2	2:B:2156:HOH:O	1.91	0.71
1:C:85:ARG:HD2	2:C:2135:HOH:O	1.91	0.70
1:C:9:SER:HB3	1:C:10:THR:HA	1.73	0.70
1:A:56:VAL:HG12	2:A:2044:HOH:O	1.91	0.69
1:D:159:PRO:HB2	1:D:161:MET:HE2	1.73	0.69
1:C:20:LEU:HD23	1:C:49:LEU:HD21	1.73	0.69
1:A:204:THR:HG21	2:A:2213:HOH:O	1.92	0.69
1:D:88:MET:HE3	1:D:100:LEU:HB3	1.77	0.66
1:C:161:MET:CE	1:C:223:ILE:HG22	2.25	0.66
1:D:96:GLU:HG3	1:D:161:MET:CE	2.28	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:203:ASN:OD1	1:A:218:HIS:HE1	1.81	0.64
1:C:9:SER:CB	1:C:10:THR:HA	2.27	0.64
1:C:161:MET:CE	2:C:2175:HOH:O	2.45	0.64
1:D:153:VAL:HB	1:D:166:THR:CG2	2.27	0.63
1:D:159:PRO:HB2	1:D:161:MET:CE	2.31	0.60
1:D:88:MET:HE3	1:D:125:THR:CG2	2.25	0.60
1:D:88:MET:HE1	1:D:101:GLN:HA	1.84	0.60
1:D:208:ALA:HB3	1:D:261:ARG:HG2	1.84	0.60
1:A:218:HIS:HD2	1:A:229:ASP:OD1	1.85	0.59
1:C:20:LEU:HD23	1:C:49:LEU:CD1	2.28	0.59
1:D:68:LEU:HD22	1:D:119:ILE:HG13	1.84	0.58
1:B:68:LEU:HD22	1:B:119:ILE:HG13	1.86	0.58
1:D:43:LEU:HD22	1:D:177:LEU:CD1	2.33	0.58
1:D:15:ARG:HH21	1:D:45:SER:HB2	1.68	0.58
1:A:15:ARG:HG2	1:A:15:ARG:O	2.04	0.57
1:B:49:LEU:HD11	1:B:267:VAL:HG22	1.86	0.57
1:C:218:HIS:HD2	1:C:229:ASP:OD1	1.87	0.57
1:A:31:GLN:HG2	2:A:2208:HOH:O	2.05	0.56
1:C:8:THR:HG22	2:C:2193:HOH:O	2.05	0.56
1:D:58:PRO:HB2	1:D:226:VAL:HG21	1.85	0.56
1:A:225:ILE:HG13	1:A:226:VAL:HG23	1.86	0.56
1:D:28:VAL:HB	1:D:36:ALA:HB3	1.87	0.55
1:C:49:LEU:HD12	1:C:267:VAL:HG22	1.86	0.55
1:A:146:GLU:HG2	1:B:201:VAL:HB	1.89	0.55
1:D:56:VAL:HG12	2:D:2027:HOH:O	2.06	0.55
1:A:208:ALA:HB3	1:A:261:ARG:HG2	1.88	0.55
1:C:161:MET:HE1	2:C:2175:HOH:O	2.07	0.55
1:A:203:ASN:OD1	1:A:218:HIS:CE1	2.58	0.55
1:B:49:LEU:HD12	1:B:267:VAL:HG22	1.89	0.55
1:C:225:ILE:HD12	1:C:225:ILE:H	1.72	0.55
1:C:9:SER:HB2	1:C:10:THR:HB	1.88	0.54
1:A:68:LEU:HD22	1:A:119:ILE:HG13	1.90	0.54
1:B:192:ARG:NH1	2:B:2231:HOH:O	2.42	0.52
1:C:11:LEU:HD13	1:C:181:GLN:NE2	2.25	0.52
1:C:203:ASN:OD1	1:C:218:HIS:HE1	1.93	0.51
1:A:15:ARG:O	1:A:15:ARG:CG	2.59	0.51
1:A:40:PHE:CD2	1:A:267:VAL:HG11	2.46	0.51
1:C:31:GLN:HB3	1:C:34:LEU:HD12	1.92	0.51
1:B:208:ALA:HB3	1:B:261:ARG:HG2	1.94	0.50
1:B:138:ALA:O	1:B:142:GLN:HG3	2.11	0.50
1:B:49:LEU:CD1	1:B:267:VAL:CG2	2.88	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:269:GLN:O	1:B:273:LYS:HG3	2.11	0.49
1:D:43:LEU:CD2	1:D:177:LEU:HD13	2.42	0.49
1:C:9:SER:CB	1:C:10:THR:CA	2.89	0.49
1:C:20:LEU:HD23	1:C:49:LEU:CD2	2.43	0.49
1:D:202:THR:O	1:D:202:THR:HG23	2.13	0.48
1:C:159:PRO:HB2	1:C:161:MET:HE3	1.95	0.48
1:C:94:ALA:HB1	1:C:159:PRO:HD3	1.96	0.48
1:D:204:THR:O	1:D:261:ARG:NE	2.46	0.47
1:D:88:MET:CE	1:D:125:THR:CG2	2.55	0.47
1:A:72:PHE:HB3	1:A:189:SER:HB3	1.97	0.47
1:C:20:LEU:HB3	1:C:49:LEU:HD21	1.97	0.47
1:C:225:ILE:HB	1:C:226:VAL:HG23	1.97	0.47
1:C:224:GLY:HA3	1:C:251:TYR:CD2	2.51	0.46
1:B:116:GLU:OE1	2:B:2160:HOH:O	2.21	0.46
1:B:72:PHE:HB3	1:B:189:SER:HB3	1.98	0.45
1:C:20:LEU:CD2	1:C:49:LEU:CD1	2.80	0.45
1:D:206:LEU:N	1:D:207:PRO:CD	2.79	0.45
1:D:188:ARG:HB2	2:D:2080:HOH:O	2.15	0.45
1:D:11:LEU:HD11	1:D:177:LEU:HD11	1.98	0.45
1:C:218:HIS:CD2	1:C:229:ASP:OD1	2.69	0.45
1:B:96:GLU:HB3	2:B:2137:HOH:O	2.18	0.44
1:C:20:LEU:HG	1:C:24:ILE:HD12	1.98	0.44
1:D:11:LEU:HA	1:D:11:LEU:HD12	1.89	0.44
1:A:25:GLN:NE2	1:A:51:VAL:HG13	2.32	0.44
1:D:56:VAL:HG23	1:D:166:THR:OG1	2.17	0.43
1:A:15:ARG:HG2	1:A:47:ALA:HA	2.00	0.43
1:C:258:GLU:O	1:C:262:GLN:HG3	2.18	0.43
1:B:49:LEU:HD12	1:B:267:VAL:CG2	2.48	0.43
1:D:152:THR:HA	1:D:166:THR:O	2.18	0.43
1:C:26:GLN:HG3	2:C:2028:HOH:O	2.19	0.43
1:C:180:GLY:HA3	2:C:2189:HOH:O	2.18	0.43
1:D:162:LYS:HB3	2:D:2067:HOH:O	2.18	0.42
1:B:68:LEU:HD23	1:B:193:LEU:HG	2.00	0.42
1:B:258:GLU:O	1:B:262:GLN:HG3	2.20	0.42
1:C:56:VAL:HG22	1:C:151:ASN:CG	2.39	0.42
1:D:159:PRO:CB	1:D:161:MET:HE2	2.47	0.42
1:B:94:ALA:HB1	1:B:159:PRO:HD3	2.02	0.42
1:C:224:GLY:HA3	1:C:251:TYR:CE2	2.54	0.42
1:D:167:THR:HG23	1:D:168:SER:N	2.34	0.42
1:C:118:MET:O	1:C:122:SER:HA	2.20	0.42
1:D:24:ILE:HD12	1:D:49:LEU:HD22	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:265:ARG:O	1:D:269:GLN:HG3	2.19	0.41
1:A:192:ARG:NH1	2:A:2171:HOH:O	2.53	0.41
1:A:99:THR:HG22	1:C:99:THR:HG22	2.03	0.41
1:A:31:GLN:HB2	1:A:31:GLN:HE21	1.66	0.41
1:D:224:GLY:HA2	1:D:252:ASN:ND2	2.35	0.41
1:C:31:GLN:NE2	2:C:2030:HOH:O	2.54	0.41
1:B:216:ILE:HG12	1:B:218:HIS:HB3	2.01	0.41
1:C:58:PRO:HG2	1:C:226:VAL:HG21	2.03	0.40
1:A:204:THR:O	1:A:204:THR:CG2	2.68	0.40
1:D:72:PHE:HB3	1:D:189:SER:HB3	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:113:GLU:OE1	1:D:1:MET:O[1_565]	2.17	0.03

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	263/298 (88%)	256 (97%)	7 (3%)	0	100	100
1	B	263/298 (88%)	256 (97%)	7 (3%)	0	100	100
1	C	268/298 (90%)	256 (96%)	9 (3%)	3 (1%)	14	5
1	D	272/298 (91%)	265 (97%)	6 (2%)	1 (0%)	34	24
All	All	1066/1192 (89%)	1033 (97%)	29 (3%)	4 (0%)	34	24

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	7	PRO
1	C	9	SER
1	D	32	PRO
1	C	223	ILE

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	222/252 (88%)	219 (99%)	3 (1%)	67	65
1	B	222/252 (88%)	219 (99%)	3 (1%)	67	65
1	C	226/252 (90%)	217 (96%)	9 (4%)	31	22
1	D	229/252 (91%)	227 (99%)	2 (1%)	78	79
All	All	899/1008 (89%)	882 (98%)	17 (2%)	57	53

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	LEU
1	A	31	GLN
1	A	122	SER
1	B	11	LEU
1	B	63	ILE
1	B	117	LEU
1	C	8	THR
1	C	10	THR
1	C	11	LEU
1	C	26	GLN
1	C	122	SER
1	C	188	ARG
1	C	193	LEU
1	C	225	ILE
1	C	252	ASN
1	D	82	LEU
1	D	122	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	GLN
1	A	31	GLN
1	A	55	GLN
1	A	218	HIS
1	A	269	GLN
1	B	19	ASN
1	B	106	ASN
1	C	33	ASN
1	C	181	GLN
1	C	218	HIS
1	C	252	ASN
1	D	108	GLN
1	D	252	ASN
1	D	269	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	265/298 (88%)	-0.26	5 (1%) 66 69	13, 25, 45, 53	0
1	B	265/298 (88%)	-0.21	3 (1%) 80 82	12, 20, 31, 51	0
1	C	270/298 (90%)	-0.13	13 (4%) 30 33	12, 21, 42, 75	0
1	D	274/298 (91%)	0.26	16 (5%) 23 25	19, 39, 59, 74	0
All	All	1074/1192 (90%)	-0.08	37 (3%) 45 48	12, 25, 50, 75	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	6	ALA	11.3
1	C	8	THR	7.7
1	D	1	MET	7.5
1	C	7	PRO	6.9
1	D	2	PRO	5.4
1	D	274	LEU	5.0
1	D	273	LYS	5.0
1	D	32	PRO	4.6
1	A	274	LEU	4.5
1	D	30	ARG	4.2
1	C	9	SER	3.9
1	C	10	THR	3.8
1	A	10	THR	3.8
1	B	275	SER	3.6
1	C	225	ILE	3.3
1	C	33	ASN	3.3
1	C	274	LEU	3.3
1	C	275	SER	3.1
1	D	13	PRO	2.9
1	A	251	TYR	2.8
1	D	27	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	32	PRO	2.7
1	D	34	LEU	2.6
1	D	33	ASN	2.5
1	D	270	ALA	2.5
1	B	274	LEU	2.5
1	A	15	ARG	2.4
1	C	26	GLN	2.4
1	D	262	GLN	2.3
1	D	26	GLN	2.2
1	D	31	GLN	2.2
1	A	273	LYS	2.2
1	C	49	LEU	2.1
1	D	23	GLN	2.1
1	B	273	LYS	2.0
1	C	251	TYR	2.0
1	D	11	LEU	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.