



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:03 PM GMT

PDB ID : 4NL8
Title : PriA Helicase Bound to SSB C-terminal Tail Peptide
Authors : Bhattacharyya, B.; George, N.P.; Thurmes, T.M.; Keck, J.L.
Deposited on : 2013-11-13
Resolution : 4.08 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
<http://wwpdb.org/validation/2016/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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

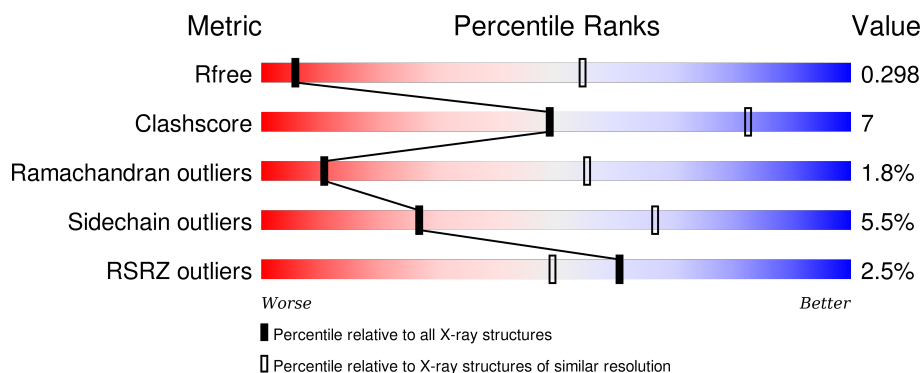
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION






The reported resolution of this entry is 4.08 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}	91344	1009 (4.52-3.60)
Clashscore	102246	1107 (4.52-3.60)
Ramachandran outliers	100387	1053 (4.52-3.60)
Sidechain outliers	100360	1039 (4.52-3.60)
RSRZ outliers	91569	1012 (4.52-3.60)

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. 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	C	10	
1	D	10	
1	F	10	
2	A	747	
2	B	747	

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Mol	Chain	Length	Quality of chain
2	E	747	<div><div></div><div>3%</div><div>63%</div><div>11%</div><div>•</div><div>24%</div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13380 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 Single-stranded DNA-binding protein.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	C	3	Total	C	N	O	0	0	0
			27	20	3	4			
1	D	5	Total	C	N	O	0	0	0
			43	28	5	10			
1	F	3	Total	C	N	O	0	0	0
			27	20	3	4			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	168	TRP	-	EXPRESSION TAG	UNP G8W6B0
D	168	TRP	-	EXPRESSION TAG	UNP G8W6B0
F	168	TRP	-	EXPRESSION TAG	UNP G8W6B0

- Molecule 2 is a protein called Primosome assembly protein PriA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	549	Total	C	N	O	S	0	1	0
			4335	2760	788	772	15			
2	B	566	Total	C	N	O	S	0	1	0
			4449	2842	805	787	15			
2	E	571	Total	C	N	O	S	0	1	0
			4493	2870	814	794	15			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-15	HIS	-	EXPRESSION TAG	UNP A6TGC5
A	-14	HIS	-	EXPRESSION TAG	UNP A6TGC5
A	-13	HIS	-	EXPRESSION TAG	UNP A6TGC5
A	-12	HIS	-	EXPRESSION TAG	UNP A6TGC5
A	-11	HIS	-	EXPRESSION TAG	UNP A6TGC5

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	HIS	-	EXPRESSION TAG	UNP A6TGC5
A	-9	SER	-	EXPRESSION TAG	UNP A6TGC5
A	-8	SER	-	EXPRESSION TAG	UNP A6TGC5
A	-7	GLY	-	EXPRESSION TAG	UNP A6TGC5
A	-6	LEU	-	EXPRESSION TAG	UNP A6TGC5
A	-5	VAL	-	EXPRESSION TAG	UNP A6TGC5
A	-4	PRO	-	EXPRESSION TAG	UNP A6TGC5
A	-3	ARG	-	EXPRESSION TAG	UNP A6TGC5
A	-2	GLY	-	EXPRESSION TAG	UNP A6TGC5
A	-1	SER	-	EXPRESSION TAG	UNP A6TGC5
A	0	HIS	-	EXPRESSION TAG	UNP A6TGC5
B	-15	HIS	-	EXPRESSION TAG	UNP A6TGC5
B	-14	HIS	-	EXPRESSION TAG	UNP A6TGC5
B	-13	HIS	-	EXPRESSION TAG	UNP A6TGC5
B	-12	HIS	-	EXPRESSION TAG	UNP A6TGC5
B	-11	HIS	-	EXPRESSION TAG	UNP A6TGC5
B	-10	HIS	-	EXPRESSION TAG	UNP A6TGC5
B	-9	SER	-	EXPRESSION TAG	UNP A6TGC5
B	-8	SER	-	EXPRESSION TAG	UNP A6TGC5
B	-7	GLY	-	EXPRESSION TAG	UNP A6TGC5
B	-6	LEU	-	EXPRESSION TAG	UNP A6TGC5
B	-5	VAL	-	EXPRESSION TAG	UNP A6TGC5
B	-4	PRO	-	EXPRESSION TAG	UNP A6TGC5
B	-3	ARG	-	EXPRESSION TAG	UNP A6TGC5
B	-2	GLY	-	EXPRESSION TAG	UNP A6TGC5
B	-1	SER	-	EXPRESSION TAG	UNP A6TGC5
B	0	HIS	-	EXPRESSION TAG	UNP A6TGC5
E	-15	HIS	-	EXPRESSION TAG	UNP A6TGC5
E	-14	HIS	-	EXPRESSION TAG	UNP A6TGC5
E	-13	HIS	-	EXPRESSION TAG	UNP A6TGC5
E	-12	HIS	-	EXPRESSION TAG	UNP A6TGC5
E	-11	HIS	-	EXPRESSION TAG	UNP A6TGC5
E	-10	HIS	-	EXPRESSION TAG	UNP A6TGC5
E	-9	SER	-	EXPRESSION TAG	UNP A6TGC5
E	-8	SER	-	EXPRESSION TAG	UNP A6TGC5
E	-7	GLY	-	EXPRESSION TAG	UNP A6TGC5
E	-6	LEU	-	EXPRESSION TAG	UNP A6TGC5
E	-5	VAL	-	EXPRESSION TAG	UNP A6TGC5
E	-4	PRO	-	EXPRESSION TAG	UNP A6TGC5
E	-3	ARG	-	EXPRESSION TAG	UNP A6TGC5
E	-2	GLY	-	EXPRESSION TAG	UNP A6TGC5
E	-1	SER	-	EXPRESSION TAG	UNP A6TGC5

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Chain	Residue	Modelled	Actual	Comment	Reference
E	0	HIS	-	EXPRESSION TAG	UNP A6TGC5

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

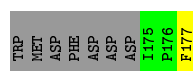
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total 2	Zn 2	0	0
3	A	2	Total 2	Zn 2	0	0
3	E	2	Total 2	Zn 2	0	0

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 errors 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.

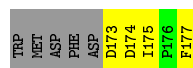
- Molecule 1: Single-stranded DNA-binding protein

Chain C: 



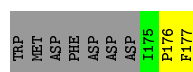
- Molecule 1: Single-stranded DNA-binding protein

Chain D: 



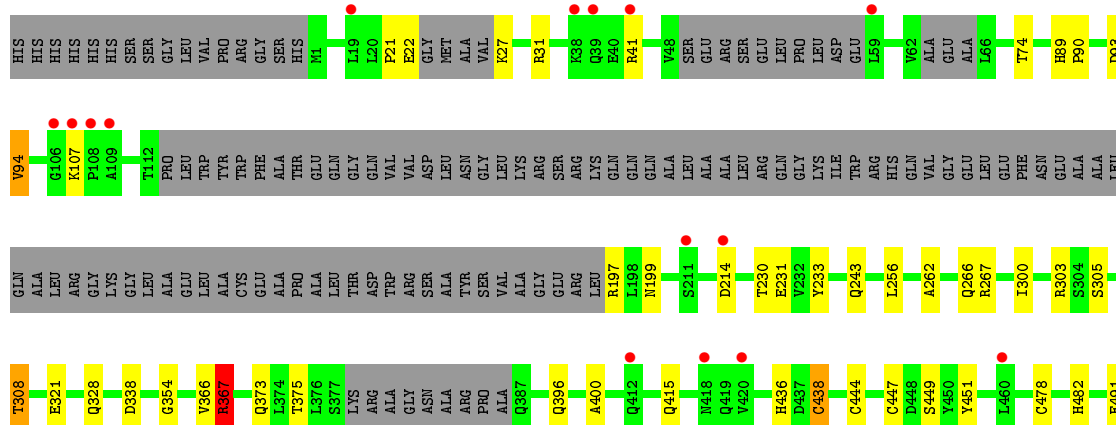
- Molecule 1: Single-stranded DNA-binding protein

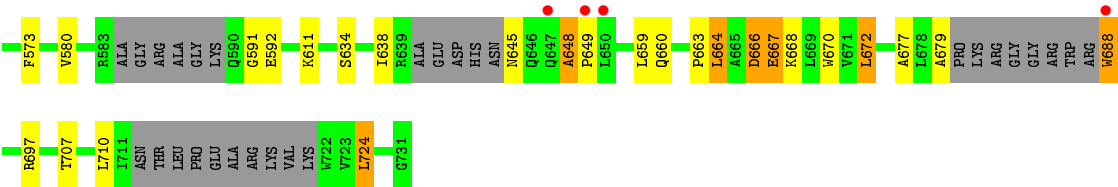
Chain F: 



- Molecule 2: Primosome assembly protein PriA

Chain A: 





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	106.61Å 153.25Å 192.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 4.08 48.71 – 4.08	Depositor EDS
% Data completeness (in resolution range)	84.2 (50.00-4.08) 84.3 (48.71-4.08)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.82 (at 4.14Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.240 , 0.298 0.249 , 0.298	Depositor DCC
R_{free} test set	1103 reflections (5.34%)	DCC
Wilson B-factor (Å ²)	124.8	Xtriage
Anisotropy	0.301	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 102.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 21752 reflections	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	13380	wwPDB-VP
Average B, all atoms (Å ²)	143.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.09% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section:
ZN

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	C	0.78	0/28	0.67	0/36
1	D	0.77	0/44	0.80	0/58
1	F	0.64	0/28	0.50	0/36
2	A	0.48	1/4439 (0.0%)	0.69	2/6035 (0.0%)
2	B	0.46	0/4557	0.68	1/6197 (0.0%)
2	E	0.45	0/4604	0.66	1/6262 (0.0%)
All	All	0.46	1/13700 (0.0%)	0.68	4/18624 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	338	ASP	CB-CG	6.75	1.66	1.51

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	338	ASP	CB-CG-OD2	6.20	123.88	118.30
2	E	367	ARG	NE-CZ-NH1	5.73	123.17	120.30
2	A	367	ARG	NE-CZ-NH1	5.58	123.09	120.30
2	B	367	ARG	NE-CZ-NH1	5.45	123.02	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	C	27	0	26	2	0
1	D	43	0	34	0	0
1	F	27	0	26	1	0
2	A	4335	0	4334	60	0
2	B	4449	0	4452	62	0
2	E	4493	0	4496	71	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	E	2	0	0	0	0
All	All	13380	0	13368	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 7.

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)
2:B:62:VAL:HG12	2:B:63:ALA:H	1.32	0.93
2:E:494:GLU:HG3	2:E:505:PRO:HG3	1.50	0.92
2:B:63:ALA:O	2:B:64:GLU:HG3	1.69	0.92
2:B:708:LEU:O	2:B:711:ILE:HD11	1.71	0.88
2:E:502:PRO:O	2:E:503:GLU:HG3	1.74	0.88
2:B:64:GLU:OE1	2:B:66:LEU:HD21	1.74	0.85
2:A:639:ARG:O	2:A:722:TRP:HA	1.75	0.85
2:A:708:LEU:O	2:A:711:ILE:HD11	1.85	0.76
2:E:526:ALA:HB1	2:E:529:ARG:HE	1.50	0.76
2:B:33:ARG:HE	2:B:64:GLU:CD	1.89	0.74
2:A:710:LEU:C	2:A:711:ILE:HG13	2.09	0.72
2:E:494:GLU:CG	2:E:505:PRO:HG3	2.19	0.72
2:A:663:PRO:O	2:A:664:LEU:HB2	1.89	0.72
2:E:498:ALA:HB1	2:E:503:GLU:O	1.90	0.71
2:B:663:PRO:O	2:B:664:LEU:HB2	1.91	0.71
2:E:387:GLN:HG2	2:E:592:GLU:HA	1.73	0.70
2:E:663:PRO:O	2:E:664:LEU:HB2	1.91	0.70
2:B:64:GLU:OE1	2:B:66:LEU:CD2	2.40	0.69
2:E:648:ALA:HB3	2:E:649:PRO:HD3	1.75	0.69
2:B:63:ALA:O	2:B:64:GLU:CG	2.41	0.68
2:E:444:CYS:O	2:E:444:CYS:SG	2.52	0.68
2:E:447:CYS:O	2:E:448:ASP:C	2.31	0.67
2:E:303:ARG:HH11	2:E:328:GLN:HE21	1.43	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:502:PRO:O	2:E:503:GLU:CG	2.42	0.67
2:B:303:ARG:HH11	2:B:328:GLN:HE21	1.43	0.67
2:B:305:SER:O	2:B:308:THR:HG23	1.94	0.67
2:A:305:SER:O	2:A:308:THR:HG23	1.95	0.66
2:E:305:SER:O	2:E:308:THR:HG23	1.96	0.66
2:E:387:GLN:HG3	2:E:592:GLU:HG3	1.78	0.66
2:A:642:ASP:CB	2:A:648:ALA:HB2	2.26	0.65
2:A:303:ARG:HH11	2:A:328:GLN:HE21	1.43	0.65
2:B:464:HIS:O	2:B:672:LEU:HD21	1.97	0.64
2:E:526:ALA:HB1	2:E:529:ARG:NE	2.11	0.64
2:A:642:ASP:HB2	2:A:648:ALA:HB2	1.79	0.63
2:E:438:CYS:HB3	2:E:478:CYS:SG	2.39	0.62
2:B:533:ARG:HA	2:B:533:ARG:HE	1.63	0.62
2:A:438:CYS:HB3	2:A:478:CYS:SG	2.40	0.61
2:A:713:THR:HG22	2:A:714:LEU:O	2.00	0.61
2:B:438:CYS:HB3	2:B:478:CYS:SG	2.40	0.61
1:C:177:PHE:OXT	1:C:177:PHE:HD1	1.84	0.61
2:E:387:GLN:CG	2:E:592:GLU:HA	2.31	0.60
2:E:386:ALA:O	2:E:387:GLN:HB2	2.01	0.60
2:A:714:LEU:HD23	2:A:714:LEU:H	1.68	0.59
2:B:533:ARG:NE	2:B:533:ARG:HA	2.19	0.58
2:A:708:LEU:O	2:A:711:ILE:CD1	2.51	0.57
2:E:555:LEU:HD23	2:E:573:PHE:CE1	2.42	0.55
2:B:710:LEU:C	2:B:711:ILE:HD13	2.27	0.55
2:B:710:LEU:O	2:B:711:ILE:HD13	2.06	0.55
1:F:176:PRO:O	1:F:177:PHE:HB2	2.06	0.55
2:E:649:PRO:HG3	2:E:688:TRP:CE3	2.41	0.55
2:A:714:LEU:N	2:A:714:LEU:HD23	2.22	0.55
1:C:177:PHE:OXT	1:C:177:PHE:CD1	2.60	0.55
2:E:464:HIS:O	2:E:672:LEU:HD21	2.07	0.55
2:B:303:ARG:NH1	2:B:328:GLN:HE21	2.05	0.54
2:B:62:VAL:CG1	2:B:63:ALA:H	2.13	0.54
2:B:582:GLY:O	2:B:584:ALA:N	2.41	0.54
2:E:303:ARG:NH1	2:E:328:GLN:HE21	2.04	0.54
2:E:387:GLN:HG2	2:E:591:GLY:O	2.08	0.53
2:A:243:GLN:HA	2:E:466:ASP:O	2.08	0.53
2:A:400:ALA:HB2	2:A:496:ALA:HB1	1.91	0.53
2:B:411:ARG:NH2	2:B:500:LEU:O	2.42	0.53
2:E:666:ASP:O	2:E:667:GLU:HG2	2.09	0.53
2:E:634:SER:OG	2:E:697:ARG:HG3	2.09	0.53
2:A:303:ARG:NH1	2:A:328:GLN:HE21	2.06	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:62:VAL:HG12	2:B:63:ALA:N	2.13	0.52
2:B:666:ASP:O	2:B:667:GLU:HG2	2.10	0.52
2:E:498:ALA:CB	2:E:503:GLU:O	2.58	0.52
2:B:634:SER:OG	2:B:697:ARG:HG3	2.09	0.52
2:B:579:GLN:HG2	2:B:583:ARG:NH2	2.24	0.52
2:A:666:ASP:O	2:A:667:GLU:HG2	2.09	0.52
2:E:411:ARG:NH2	2:E:500:LEU:O	2.43	0.52
2:B:555:LEU:HD23	2:B:573:PHE:CE1	2.45	0.51
2:B:21:PRO:O	2:B:22:GLU:O	2.28	0.51
2:A:634:SER:OG	2:A:697:ARG:HG3	2.10	0.51
2:B:447:CYS:SG	2:B:465:CYS:HB3	2.51	0.51
2:B:400:ALA:HB2	2:B:496:ALA:HB1	1.93	0.51
2:B:33:ARG:NE	2:B:64:GLU:OE1	2.43	0.51
2:E:386:ALA:O	2:E:387:GLN:CB	2.58	0.51
2:B:677:ALA:C	2:B:679:ALA:H	2.14	0.50
2:E:502:PRO:O	2:E:503:GLU:CB	2.59	0.50
2:E:501:PHE:O	2:E:502:PRO:C	2.48	0.50
2:A:645:ASN:ND2	2:A:645:ASN:C	2.65	0.50
2:A:666:ASP:C	2:A:668:LYS:H	2.14	0.50
2:E:666:ASP:C	2:E:668:LYS:H	2.15	0.50
2:E:447:CYS:SG	2:E:465:CYS:HB3	2.51	0.49
2:A:646:GLN:O	2:A:648:ALA:N	2.41	0.49
2:E:400:ALA:HB2	2:E:496:ALA:HB1	1.93	0.49
2:E:233:TYR:OH	2:E:354:GLY:HA3	2.13	0.48
2:E:491:GLU:O	2:E:492:GLN:C	2.51	0.48
2:B:638:ILE:HG12	2:B:724:LEU:HD22	1.95	0.48
2:B:666:ASP:C	2:B:668:LYS:H	2.16	0.48
2:E:400:ALA:HB2	2:E:496:ALA:CB	2.43	0.48
2:B:533:ARG:HD3	2:B:534:ILE:H	1.79	0.48
2:B:491:GLU:O	2:B:492:GLN:C	2.52	0.48
2:A:710:LEU:O	2:A:711:ILE:HG13	2.13	0.48
2:E:447:CYS:O	2:E:449:SER:N	2.46	0.48
2:A:712:ASN:O	2:A:713:THR:OG1	2.24	0.48
2:E:659:LEU:HD21	2:E:707:THR:HG21	1.96	0.47
2:B:646:GLN:O	2:B:648:ALA:N	2.42	0.47
2:A:400:ALA:HB2	2:A:496:ALA:CB	2.45	0.47
2:B:400:ALA:HB2	2:B:496:ALA:CB	2.44	0.47
2:E:504:VAL:HB	2:E:505:PRO:CD	2.45	0.47
2:A:659:LEU:HD21	2:A:707:THR:HG21	1.97	0.47
2:A:233:TYR:OH	2:A:354:GLY:HA3	2.14	0.47
2:A:677:ALA:C	2:A:679:ALA:H	2.18	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:491:GLU:O	2:A:492:GLN:C	2.52	0.47
2:E:447:CYS:O	2:E:449:SER:CB	2.64	0.46
2:E:638:ILE:HG12	2:E:724:LEU:HD22	1.96	0.46
2:A:648:ALA:O	2:A:651:PHE:HB3	2.16	0.46
2:A:648:ALA:HB3	2:A:649:PRO:HD3	1.97	0.46
2:A:243:GLN:HA	2:E:467:SER:HA	1.97	0.46
2:E:367:ARG:HH11	2:E:367:ARG:CB	2.28	0.46
2:A:646:GLN:HB3	2:A:647:GLN:HE21	1.81	0.46
2:B:367:ARG:HH11	2:B:367:ARG:CB	2.29	0.46
2:E:504:VAL:HB	2:E:505:PRO:HD2	1.98	0.46
2:B:646:GLN:HB3	2:B:647:GLN:HE21	1.80	0.46
2:B:93:ASP:O	2:B:94:VAL:C	2.54	0.46
2:A:266:GLN:NE2	2:B:441:ILE:HD12	2.31	0.46
2:B:659:LEU:HD21	2:B:707:THR:HG21	1.97	0.46
2:E:677:ALA:C	2:E:679:ALA:H	2.20	0.46
2:A:498:ALA:N	2:A:499:PRO:HD2	2.30	0.46
2:A:651:PHE:HE2	2:A:714:LEU:C	2.19	0.46
2:A:367:ARG:CB	2:A:367:ARG:HH11	2.30	0.45
2:B:31:ARG:NH2	2:B:107:LYS:O	2.44	0.45
2:E:254:ILE:HD12	2:E:542:ALA:HB2	1.99	0.45
2:E:387:GLN:HG2	2:E:592:GLU:CA	2.44	0.45
2:A:638:ILE:HG12	2:A:724:LEU:HD22	1.98	0.45
2:A:679:ALA:HB3	2:A:687:ARG:HD3	1.99	0.45
2:A:670:TRP:CZ3	2:A:672:LEU:HD23	2.51	0.45
2:E:367:ARG:HB2	2:E:367:ARG:HH11	1.82	0.45
2:A:93:ASP:O	2:A:94:VAL:C	2.55	0.45
2:B:498:ALA:N	2:B:499:PRO:HD2	2.33	0.44
2:B:648:ALA:HB3	2:B:649:PRO:HD3	1.99	0.44
2:E:21:PRO:O	2:E:22:GLU:C	2.55	0.44
2:E:444:CYS:SG	2:E:449:SER:HB3	2.58	0.44
2:B:233:TYR:OH	2:B:354:GLY:HA3	2.16	0.44
2:E:93:ASP:O	2:E:94:VAL:C	2.55	0.44
2:B:367:ARG:HH11	2:B:367:ARG:HB2	1.83	0.44
2:B:197:ARG:HG2	2:B:197:ARG:H	1.54	0.44
2:E:31:ARG:NH2	2:E:107:LYS:O	2.44	0.43
2:B:254:ILE:HD12	2:B:542:ALA:HB2	2.00	0.43
2:A:214[B]:ASP:OD2	2:E:457:GLN:HG2	2.17	0.43
2:E:649:PRO:HG3	2:E:688:TRP:CD2	2.53	0.43
2:E:498:ALA:N	2:E:499:PRO:HD2	2.33	0.43
2:B:670:TRP:CZ3	2:B:672:LEU:HD23	2.53	0.43
2:E:21:PRO:O	2:E:22:GLU:O	2.37	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:367:ARG:HB2	2:A:367:ARG:HH11	1.84	0.43
2:B:475:CYS:HB2	2:B:483:LEU:HD21	2.01	0.43
2:A:651:PHE:CE2	2:A:714:LEU:C	2.91	0.43
2:E:506:ILE:HG22	2:E:507:SER:N	2.34	0.43
2:E:231:GLU:OE2	2:E:267:ARG:NH2	2.52	0.43
2:A:444:CYS:HA	2:A:451:TYR:CE2	2.54	0.43
2:E:444:CYS:HA	2:E:451:TYR:CE2	2.54	0.43
2:B:89:HIS:CG	2:B:90:PRO:HD2	2.53	0.43
2:A:256:LEU:HD21	2:A:300:ILE:HG22	2.01	0.42
2:A:231:GLU:OE2	2:A:267:ARG:NH2	2.51	0.42
2:A:305:SER:O	2:A:308:THR:CG2	2.65	0.42
2:A:366:VAL:HG11	2:A:373:GLN:HE21	1.84	0.42
2:A:447:CYS:SG	2:A:449:SER:HB2	2.60	0.42
2:E:256:LEU:HD21	2:E:300:ILE:HG22	2.01	0.42
2:B:648:ALA:O	2:B:651:PHE:HB3	2.19	0.42
2:A:21:PRO:O	2:A:22:GLU:C	2.58	0.42
2:A:89:HIS:CG	2:A:90:PRO:HD2	2.55	0.42
2:B:33:ARG:O	2:B:63:ALA:HB3	2.20	0.42
2:E:305:SER:O	2:E:308:THR:CG2	2.66	0.41
2:B:581:SER:O	2:B:583:ARG:N	2.53	0.41
2:B:256:LEU:HD21	2:B:300:ILE:HG22	2.02	0.41
2:B:63:ALA:C	2:B:64:GLU:HG3	2.36	0.41
2:B:305:SER:O	2:B:308:THR:CG2	2.64	0.41
2:A:645:ASN:HD22	2:A:645:ASN:C	2.23	0.41
2:A:396:GLN:O	2:A:436:HIS:NE2	2.45	0.41
2:B:583:ARG:O	2:B:584:ALA:HB2	2.20	0.41
2:E:89:HIS:CG	2:E:90:PRO:HD2	2.55	0.41
2:B:366:VAL:HG11	2:B:373:GLN:HE21	1.85	0.41
2:B:231:GLU:OE2	2:B:267:ARG:NH2	2.53	0.41
2:A:720:VAL:HG12	2:A:720:VAL:O	2.20	0.41
2:E:494:GLU:HA	2:E:497:LEU:HB2	2.01	0.41
2:A:647:GLN:O	2:A:650:LEU:HB2	2.20	0.41
2:A:712:ASN:O	2:A:713:THR:CB	2.68	0.41
2:E:670:TRP:CZ3	2:E:672:LEU:HD23	2.55	0.41
2:E:659:LEU:CD2	2:E:707:THR:HG21	2.51	0.41
2:E:475:CYS:HB2	2:E:483:LEU:HD21	2.03	0.41
2:E:396:GLN:O	2:E:436:HIS:NE2	2.48	0.41
2:E:262:ALA:O	2:E:266:GLN:HG2	2.21	0.41
2:E:366:VAL:HG11	2:E:373:GLN:HE21	1.85	0.41
2:A:31:ARG:NH2	2:A:107:LYS:O	2.44	0.41
2:B:444:CYS:HA	2:B:451:TYR:CE2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:659:LEU:CD2	2:A:707:THR:HG21	2.52	0.40
2:A:21:PRO:O	2:A:22:GLU:O	2.39	0.40
2:B:494:GLU:HA	2:B:497:LEU:HB2	2.01	0.40
2:E:447:CYS:O	2:E:449:SER:HB2	2.22	0.40
2:A:262:ALA:O	2:A:266:GLN:HG2	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	C	1/10 (10%)	1 (100%)	0	0	100	100
1	D	3/10 (30%)	0	1 (33%)	2 (67%)	0	0
1	F	1/10 (10%)	0	1 (100%)	0	100	100
2	A	530/747 (71%)	493 (93%)	29 (6%)	8 (2%)	13	58
2	B	545/747 (73%)	504 (92%)	32 (6%)	9 (2%)	11	57
2	E	550/747 (74%)	512 (93%)	27 (5%)	11 (2%)	9	54
All	All	1630/2271 (72%)	1510 (93%)	90 (6%)	30 (2%)	11	55

All (30) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	664	LEU
2	A	667	GLU
2	A	713	THR
1	D	174	ASP
2	B	62	VAL
2	B	583	ARG
2	B	584	ALA

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Mol	Chain	Res	Type
2	B	664	LEU
2	B	667	GLU
2	E	387	GLN
2	E	503	GLU
2	E	664	LEU
2	E	667	GLU
2	E	448	ASP
2	E	505	PRO
2	E	527	VAL
2	E	648	ALA
2	A	492	GLN
2	A	722	TRP
2	B	492	GLN
2	B	647	GLN
2	E	492	GLN
2	A	647	GLN
1	D	175	ILE
2	B	321	GLU
2	E	502	PRO
2	A	321	GLU
2	A	94	VAL
2	E	94	VAL
2	B	94	VAL

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	C	3/10 (30%)	3 (100%)	0	100	100
1	D	5/10 (50%)	3 (60%)	2 (40%)	0	0
1	F	3/10 (30%)	3 (100%)	0	100	100
2	A	464/614 (76%)	437 (94%)	27 (6%)	25	65
2	B	473/614 (77%)	448 (95%)	25 (5%)	28	67
2	E	480/614 (78%)	456 (95%)	24 (5%)	30	69

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1428/1872 (76%)	1350 (94%)	78 (6%)	27 67

All (78) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	27	LYS
2	A	41	ARG
2	A	74	THR
2	A	197	ARG
2	A	199	ASN
2	A	230	THR
2	A	308	THR
2	A	367	ARG
2	A	375	THR
2	A	415	GLN
2	A	438	CYS
2	A	482	HIS
2	A	494	GLU
2	A	556	LEU
2	A	580	VAL
2	A	611	LYS
2	A	645	ASN
2	A	646	GLN
2	A	660	GLN
2	A	666	ASP
2	A	672	LEU
2	A	688	TRP
2	A	710	LEU
2	A	711	ILE
2	A	714	LEU
2	A	722	TRP
2	A	724	LEU
1	D	173	ASP
1	D	177	PHE
2	B	41	ARG
2	B	74	THR
2	B	197	ARG
2	B	199	ASN
2	B	230	THR
2	B	308	THR
2	B	367	ARG
2	B	375	THR

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Mol	Chain	Res	Type
2	B	415	GLN
2	B	438	CYS
2	B	447	CYS
2	B	482	HIS
2	B	494	GLU
2	B	533	ARG
2	B	556	LEU
2	B	580	VAL
2	B	611	LYS
2	B	646	GLN
2	B	660	GLN
2	B	666	ASP
2	B	672	LEU
2	B	678	LEU
2	B	710	LEU
2	B	711	ILE
2	B	724	LEU
2	E	41	ARG
2	E	74	THR
2	E	197	ARG
2	E	199	ASN
2	E	230	THR
2	E	308	THR
2	E	367	ARG
2	E	375	THR
2	E	415	GLN
2	E	438	CYS
2	E	444	CYS
2	E	447	CYS
2	E	482	HIS
2	E	494	GLU
2	E	556	LEU
2	E	580	VAL
2	E	611	LYS
2	E	645	ASN
2	E	660	GLN
2	E	666	ASP
2	E	672	LEU
2	E	688	TRP
2	E	710	LEU
2	E	724	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such

sidechains are listed below:

Mol	Chain	Res	Type
2	A	199	ASN
2	A	243	GLN
2	A	266	GLN
2	A	328	GLN
2	A	329	GLN
2	A	373	GLN
2	A	495	GLN
2	A	645	ASN
2	A	646	GLN
2	A	647	GLN
2	A	660	GLN
2	B	199	ASN
2	B	328	GLN
2	B	329	GLN
2	B	373	GLN
2	B	419	GLN
2	B	495	GLN
2	B	647	GLN
2	B	657	ASN
2	B	660	GLN
2	E	199	ASN
2	E	328	GLN
2	E	329	GLN
2	E	373	GLN
2	E	495	GLN
2	E	646	GLN
2	E	660	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

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	C	3/10 (30%)	-0.20	0 100 100	119, 119, 160, 167	0
1	D	5/10 (50%)	0.45	0 100 100	169, 192, 198, 199	0
1	F	3/10 (30%)	-0.55	0 100 100	148, 148, 187, 208	0
2	A	549/747 (73%)	-0.01	15 (2%) 58 47	80, 129, 197, 257	0
2	B	566/747 (75%)	-0.10	2 (0%) 93 90	81, 134, 206, 266	0
2	E	571/747 (76%)	0.14	25 (4%) 38 28	92, 146, 224, 285	0
All	All	1697/2271 (74%)	0.01	42 (2%) 61 49	80, 137, 210, 285	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	59	LEU	5.5
2	E	506	ILE	5.2
2	E	38	LYS	5.1
2	E	505	PRO	5.1
2	E	507	SER	5.0
2	E	8	LEU	4.4
2	E	7	ALA	4.0
2	E	37	GLY	3.8
2	E	42	ILE	3.8
2	A	107	LYS	3.5
2	A	38	LYS	3.5
2	A	39	GLN	3.5
2	E	647	GLN	3.4
2	A	106	GLY	3.4
2	E	61	PRO	3.3
2	A	41	ARG	3.3
2	E	649	PRO	3.2
2	A	108	PRO	3.0
2	E	36	PHE	2.9

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Mol	Chain	Res	Type	RSRZ
2	E	43	GLY	2.8
2	E	4	ALA	2.7
2	A	109	ALA	2.6
2	A	418	ASN	2.6
2	E	688	TRP	2.5
2	E	16	PHE	2.5
2	E	6	VAL	2.5
2	A	19	LEU	2.5
2	E	428	GLY	2.4
2	E	33	ARG	2.4
2	E	41	ARG	2.4
2	E	62	VAL	2.4
2	A	460	LEU	2.4
2	E	17	ASP	2.3
2	A	420	VAL	2.1
2	B	497	LEU	2.1
2	E	650	LEU	2.1
2	B	41	ARG	2.1
2	A	214[A]	ASP	2.1
2	A	412	GLN	2.1
2	E	202	GLN	2.0
2	E	40	GLU	2.0
2	A	211	SER	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	ZN	E	802	1/1	0.99	0.15	-0.47	102,102,102,102	0
3	ZN	A	801	1/1	0.99	0.13	-0.60	103,103,103,103	0
3	ZN	E	801	1/1	0.99	0.16	-0.76	117,117,117,117	0
3	ZN	B	802	1/1	0.98	0.12	-1.07	118,118,118,118	0
3	ZN	B	801	1/1	1.00	0.11	-1.20	100,100,100,100	0
3	ZN	A	802	1/1	0.99	0.07	-1.83	129,129,129,129	0

6.5 Other polymers [i](#)

There are no such residues in this entry.