



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 08:30 pm BST

PDB ID : 3QKR
Title : Mre11 Rad50 binding domain bound to Rad50
Authors : Williams, G.J.; Williams, R.S.; Arvai, A.; Moncalian, G.; Tainer, J.A.
Deposited on : 2011-02-01
Resolution : 3.40 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
Mogul : 1.8.5 (274361), CSD as541be (2020)
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

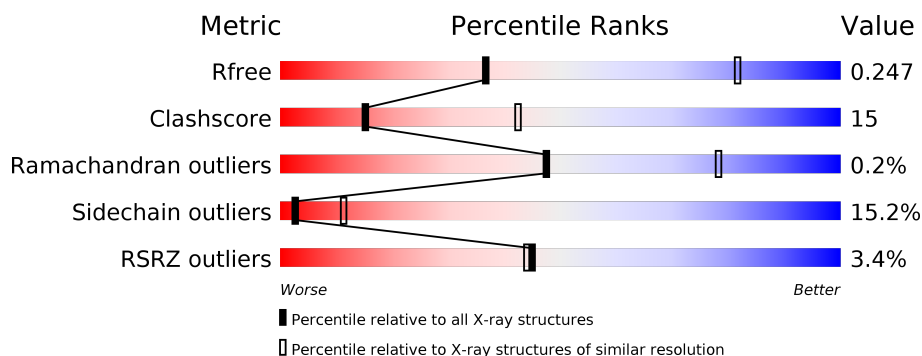
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1026 (3.48-3.32)
Clashscore	141614	1055 (3.48-3.32)
Ramachandran outliers	138981	1038 (3.48-3.32)
Sidechain outliers	138945	1038 (3.48-3.32)
RSRZ outliers	127900	2173 (3.50-3.30)

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	203	<div> <div>4%</div> <div> <div></div> <div>65%</div> <div>27%</div> <div>6%</div> </div> </div>
2	B	179	<div> <div>3%</div> <div> <div></div> <div>61%</div> <div>31%</div> <div>8%</div> </div> </div>
3	C	34	<div> <div></div> <div> <div></div> <div>50%</div> <div>35%</div> <div>9%</div> <div>6%</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3356 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 DNA double-strand break repair rad50 ATPase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	199	Total	C	N	O	S	0	0	0
			1624	1037	287	297	3			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	196	GLY	-	EXPRESSION TAG	UNP P58301
A	197	GLY	-	EXPRESSION TAG	UNP P58301
A	198	HIS	-	EXPRESSION TAG	UNP P58301
A	199	HIS	-	EXPRESSION TAG	UNP P58301
A	200	HIS	-	EXPRESSION TAG	UNP P58301
A	201	HIS	-	EXPRESSION TAG	UNP P58301
A	202	HIS	-	EXPRESSION TAG	UNP P58301
A	203	HIS	-	EXPRESSION TAG	UNP P58301

- Molecule 2 is a protein called DNA double-strand break repair rad50 ATPase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	179	Total	C	N	O	S	0	0	0
			1445	923	246	274	2			

- Molecule 3 is a protein called DNA double-strand break repair protein mre11.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	32	Total	C	N	O	0	0	0
			270	180	34	56			

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		

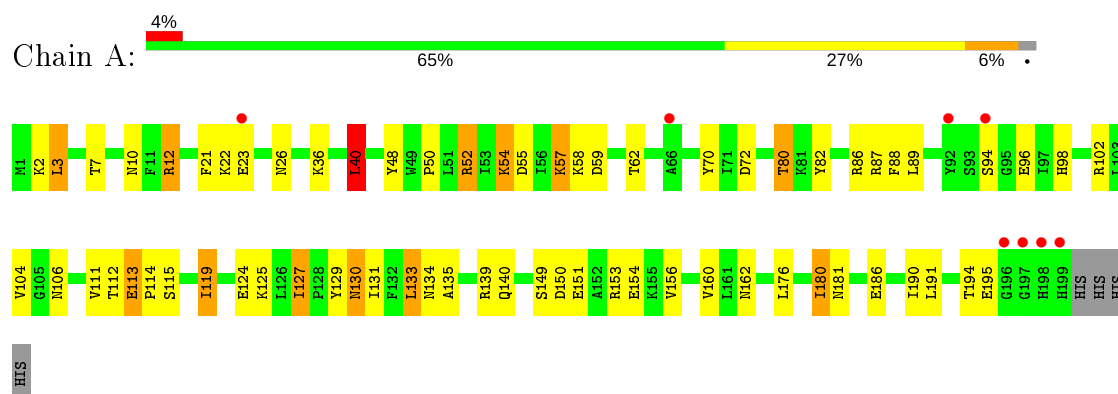
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	3	Total	O	0	0
			3	3		
5	B	4	Total	O	0	0
			4	4		

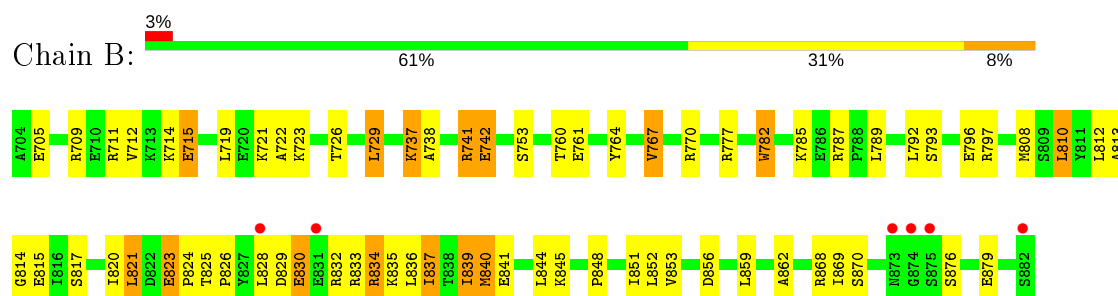
3 Residue-property plots [i](#)

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.

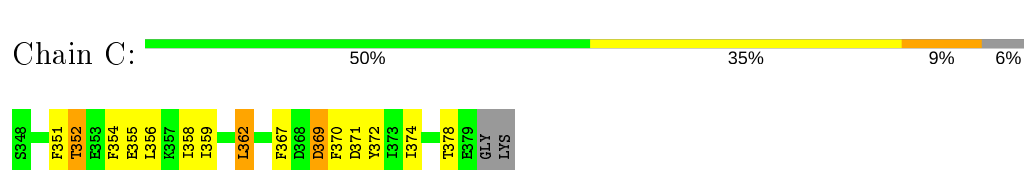
- Molecule 1: DNA double-strand break repair rad50 ATPase



- Molecule 2: DNA double-strand break repair rad50 ATPase



- Molecule 3: DNA double-strand break repair protein mre11



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	116.04Å 116.04Å 109.85Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 3.40 19.94 – 3.40	Depositor EDS
% Data completeness (in resolution range)	91.6 (50.00-3.40) 92.0 (19.94-3.40)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.46 (at 3.36Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.211 , 0.274 0.188 , 0.247	Depositor DCC
R_{free} test set	514 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å ²)	108.1	Xtriage
Anisotropy	0.209	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 99.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.045 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3356	wwPDB-VP
Average B, all atoms (Å ²)	99.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.17% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PO4

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	0.32	0/1651	0.54	1/2213 (0.0%)
2	B	0.32	0/1463	0.54	0/1961
3	C	0.38	0/275	0.61	0/370
All	All	0.33	0/3389	0.54	1/4544 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	40	LEU	CA-CB-CG	6.06	129.24	115.30

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	1624	0	1677	47	0
2	B	1445	0	1506	58	0
3	C	270	0	250	10	0
4	A	10	0	0	0	0
5	A	3	0	0	0	0
5	B	4	0	0	0	0
All	All	3356	0	3433	102	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 15.

All (102) 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:A:12:ARG:HG3	1:A:12:ARG:HH21	1.24	1.03
1:A:52:ARG:HD3	1:A:52:ARG:H	1.31	0.95
2:B:834:ARG:HB2	2:B:834:ARG:HH11	1.35	0.92
2:B:741:ARG:HG2	2:B:741:ARG:HH11	1.34	0.90
2:B:760:THR:HG21	2:B:764:TYR:HD1	1.35	0.89
1:A:89:LEU:HB2	1:A:94:SER:O	1.72	0.88
1:A:52:ARG:N	1:A:52:ARG:HD3	1.89	0.86
2:B:823:GLU:HB2	2:B:824:PRO:C	1.96	0.84
1:A:96:GLU:HG2	1:A:98:HIS:NE2	2.00	0.76
2:B:824:PRO:HG3	2:B:852:LEU:HD11	1.69	0.73
1:A:12:ARG:HH21	1:A:12:ARG:CG	2.02	0.71
2:B:841:GLU:O	2:B:845:LYS:HG2	1.90	0.71
1:A:48:TYR:OH	1:A:127:ILE:HG12	1.92	0.69
2:B:741:ARG:NH1	2:B:741:ARG:HG2	2.03	0.67
1:A:86:ARG:HB3	1:A:88:PHE:HE1	1.60	0.66
1:A:135:ALA:HB1	2:B:813:ALA:HB2	1.77	0.66
2:B:760:THR:HG21	2:B:764:TYR:CD1	2.26	0.65
1:A:180:ILE:HG23	3:C:374:ILE:HD12	1.79	0.64
2:B:770:ARG:HB3	2:B:777:ARG:HB2	1.79	0.64
1:A:12:ARG:HG3	1:A:12:ARG:NH2	2.04	0.64
3:C:351:PHE:HB3	3:C:355:GLU:HG2	1.82	0.62
2:B:830:GLU:HB3	2:B:834:ARG:NH1	2.16	0.61
1:A:129:TYR:O	1:A:133:LEU:HB2	2.01	0.61
1:A:7:THR:HB	1:A:72:ASP:HB3	1.83	0.61
1:A:52:ARG:N	1:A:52:ARG:CD	2.64	0.60
2:B:833:ARG:O	2:B:837:ILE:HG22	2.02	0.60
2:B:729:LEU:HD21	3:C:355:GLU:HG3	1.83	0.60
1:A:150:ASP:HA	1:A:153:ARG:NH2	2.16	0.60
2:B:856:ASP:O	2:B:859:LEU:HD12	2.02	0.60
2:B:737:LYS:HG3	2:B:738:ALA:N	2.17	0.59
1:A:98:HIS:O	1:A:114:PRO:HB3	2.03	0.59
1:A:130:ASN:HD22	1:A:131:ILE:H	1.49	0.59
2:B:785:LYS:HB3	2:B:787:ARG:NH2	2.18	0.58
2:B:782:TRP:O	2:B:787:ARG:NH2	2.35	0.58
1:A:180:ILE:HD12	2:B:723:LYS:HA	1.86	0.57
1:A:135:ALA:CB	2:B:813:ALA:HB2	2.34	0.56
2:B:721:LYS:HD3	3:C:378:THR:HG23	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:127:ILE:HG12	1:A:127:ILE:O	2.06	0.56
2:B:823:GLU:O	2:B:853:VAL:O	2.26	0.54
2:B:753:SER:HA	2:B:767:VAL:HG13	1.88	0.54
1:A:36:LYS:HB3	2:B:869:ILE:HG21	1.89	0.54
2:B:837:ILE:HD12	2:B:862:ALA:HB2	1.88	0.54
2:B:820:ILE:HG12	2:B:851:ILE:HB	1.90	0.54
2:B:840:MET:HA	2:B:844:LEU:HG	1.90	0.53
2:B:834:ARG:HB2	2:B:834:ARG:NH1	2.14	0.53
1:A:3:LEU:HD23	1:A:3:LEU:H	1.73	0.53
1:A:186:GLU:O	1:A:190:ILE:HD13	2.08	0.53
1:A:50:PRO:O	1:A:58:LYS:HE2	2.09	0.53
2:B:808:MET:O	2:B:812:LEU:HD12	2.09	0.53
2:B:823:GLU:HB2	2:B:824:PRO:CA	2.39	0.52
2:B:764:TYR:OH	2:B:796:GLU:OE2	2.27	0.52
2:B:789:LEU:H	2:B:789:LEU:HD12	1.74	0.52
1:A:12:ARG:NH2	1:A:12:ARG:CG	2.68	0.51
1:A:195:GLU:HB2	2:B:709:ARG:NH1	2.26	0.51
1:A:191:LEU:O	2:B:709:ARG:NH1	2.43	0.50
2:B:836:LEU:HA	2:B:839:ILE:HD13	1.94	0.50
1:A:57:LYS:HA	1:A:57:LYS:HE2	1.94	0.49
1:A:54:LYS:HG3	1:A:55:ASP:H	1.77	0.49
2:B:738:ALA:O	2:B:742:GLU:OE2	2.31	0.48
1:A:130:ASN:ND2	1:A:130:ASN:H	2.12	0.48
2:B:711:ARG:NH2	2:B:715:GLU:OE2	2.46	0.48
3:C:356:LEU:HA	3:C:359:ILE:HD12	1.96	0.48
2:B:722:ALA:O	2:B:726:THR:HG23	2.13	0.48
1:A:115:SER:O	1:A:119:ILE:HB	2.13	0.47
2:B:824:PRO:CG	2:B:852:LEU:HD11	2.41	0.47
1:A:130:ASN:ND2	1:A:130:ASN:N	2.61	0.47
1:A:134:ASN:HB3	1:A:160:VAL:HG13	1.96	0.47
1:A:191:LEU:HD21	2:B:712:VAL:HG12	1.96	0.47
2:B:742:GLU:OE2	2:B:742:GLU:N	2.39	0.46
1:A:26:ASN:N	1:A:26:ASN:HD22	2.12	0.46
2:B:742:GLU:CD	2:B:742:GLU:H	2.12	0.46
2:B:810:LEU:O	2:B:814:GLY:N	2.44	0.46
2:B:868:ARG:HB2	2:B:879:GLU:HB3	1.97	0.45
1:A:149:SER:OG	1:A:151:GLU:OE1	2.34	0.45
2:B:835:LYS:O	2:B:839:ILE:HG23	2.16	0.45
2:B:836:LEU:O	2:B:840:MET:HB3	2.17	0.45
1:A:111:VAL:HG23	1:A:112:THR:HG23	1.98	0.45
1:A:150:ASP:HA	1:A:153:ARG:HH22	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:729:LEU:CD2	3:C:355:GLU:HG3	2.47	0.45
1:A:40:LEU:HD12	2:B:853:VAL:HG21	1.99	0.44
2:B:793:SER:HB3	2:B:796:GLU:OE2	2.17	0.44
1:A:3:LEU:O	1:A:21:PHE:HB2	2.18	0.44
1:A:80:THR:HB	1:A:82:TYR:CE1	2.53	0.44
1:A:113:GLU:HA	1:A:114:PRO:HD3	1.86	0.44
1:A:48:TYR:OH	1:A:127:ILE:CG1	2.64	0.43
3:C:369:ASP:OD2	3:C:372:TYR:HB3	2.17	0.43
3:C:352:THR:OG1	3:C:355:GLU:HB3	2.19	0.43
2:B:821:LEU:HB2	2:B:852:LEU:HD12	2.00	0.43
1:A:139:ARG:HG3	1:A:140:GLN:N	2.35	0.42
1:A:70:TYR:CE1	1:A:72:ASP:HB2	2.54	0.42
3:C:358:ILE:O	3:C:362:LEU:HB2	2.19	0.42
2:B:721:LYS:HG2	2:B:722:ALA:N	2.34	0.42
2:B:870:SER:O	2:B:876:SER:HA	2.20	0.42
2:B:760:THR:O	2:B:761:GLU:HB2	2.20	0.41
2:B:792:LEU:O	2:B:797:ARG:NH2	2.53	0.41
1:A:2:LYS:HE2	1:A:23:GLU:HG3	2.03	0.41
1:A:176:LEU:HD12	3:C:370:PHE:HB3	2.03	0.41
2:B:830:GLU:O	2:B:834:ARG:N	2.45	0.40
2:B:836:LEU:HA	2:B:839:ILE:CD1	2.51	0.40
2:B:825:THR:HA	2:B:826:PRO:HD3	1.94	0.40
2:B:830:GLU:HB3	2:B:834:ARG:HH12	1.84	0.40
2:B:817:SER:HA	2:B:848:PRO:HG2	2.04	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	197/203 (97%)	186 (94%)	11 (6%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	177/179 (99%)	170 (96%)	6 (3%)	1 (1%)	25	57
3	C	30/34 (88%)	28 (93%)	2 (7%)	0	100	100
All	All	404/416 (97%)	384 (95%)	19 (5%)	1 (0%)	47	78

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	823	GLU

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	176/180 (98%)	148 (84%)	28 (16%)	2	10
2	B	156/156 (100%)	135 (86%)	21 (14%)	4	14
3	C	30/32 (94%)	24 (80%)	6 (20%)	1	3
All	All	362/368 (98%)	307 (85%)	55 (15%)	3	11

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	10	ASN
1	A	12	ARG
1	A	22	LYS
1	A	40	LEU
1	A	52	ARG
1	A	54	LYS
1	A	57	LYS
1	A	59	ASP
1	A	62	THR
1	A	80	THR
1	A	87	ARG
1	A	102	ARG

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Mol	Chain	Res	Type
1	A	104	VAL
1	A	106	ASN
1	A	113	GLU
1	A	119	ILE
1	A	124	GLU
1	A	125	LYS
1	A	127	ILE
1	A	130	ASN
1	A	133	LEU
1	A	154	GLU
1	A	156	VAL
1	A	162	ASN
1	A	180	ILE
1	A	181	ASN
1	A	194	THR
2	B	705	GLU
2	B	714	LYS
2	B	715	GLU
2	B	719	LEU
2	B	729	LEU
2	B	737	LYS
2	B	741	ARG
2	B	742	GLU
2	B	767	VAL
2	B	782	TRP
2	B	810	LEU
2	B	815	GLU
2	B	821	LEU
2	B	828	LEU
2	B	829	ASP
2	B	830	GLU
2	B	832	ARG
2	B	834	ARG
2	B	837	ILE
2	B	839	ILE
2	B	840	MET
3	C	352	THR
3	C	354	PHE
3	C	362	LEU
3	C	367	PHE
3	C	369	ASP
3	C	371	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	130	ASN
1	A	134	ASN
2	B	774	ASN
2	B	855	HIS

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](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	PO4	A	200	-	4,4,4	0.88	0	6,6,6	0.51	0
4	PO4	A	201	-	4,4,4	0.89	0	6,6,6	0.48	0

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	A	199/203 (98%)	-0.20	8 (4%)	38	37	33, 94, 173, 229	0
2	B	179/179 (100%)	-0.36	6 (3%)	45	44	28, 89, 156, 199	0
3	C	32/34 (94%)	-0.43	0	100	100	61, 106, 162, 166	0
All	All	410/416 (98%)	-0.29	14 (3%)	45	44	28, 95, 163, 229	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	199	HIS	6.1
1	A	197	GLY	6.1
1	A	92	TYR	5.5
1	A	198	HIS	5.2
2	B	873	ASN	4.1
2	B	828	LEU	3.3
2	B	882	SER	2.4
1	A	94	SER	2.4
2	B	874	GLY	2.3
2	B	831	GLU	2.2
1	A	196	GLY	2.2
1	A	66	ALA	2.2
2	B	875	SER	2.2
1	A	23	GLU	2.0

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

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

6.3 Carbohydrates [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. 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	B-factors(\AA^2)	Q<0.9
4	PO4	A	201	5/5	0.83	0.31	93,121,168,180	0
4	PO4	A	200	5/5	0.96	0.08	35,57,75,136	0

6.5 Other polymers [i](#)

There are no such residues in this entry.