



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:51 PM UTC

PDB ID : 3BX2 / pdb_00003bx2
Title : Puf4 RNA binding domain bound to HO endonuclease RNA 3' UTR recognition sequence
Authors : Miller, M.T.; Higgin, J.J.; Hall, T.M.T.
Deposited on : 2008-01-11
Resolution : 2.84 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

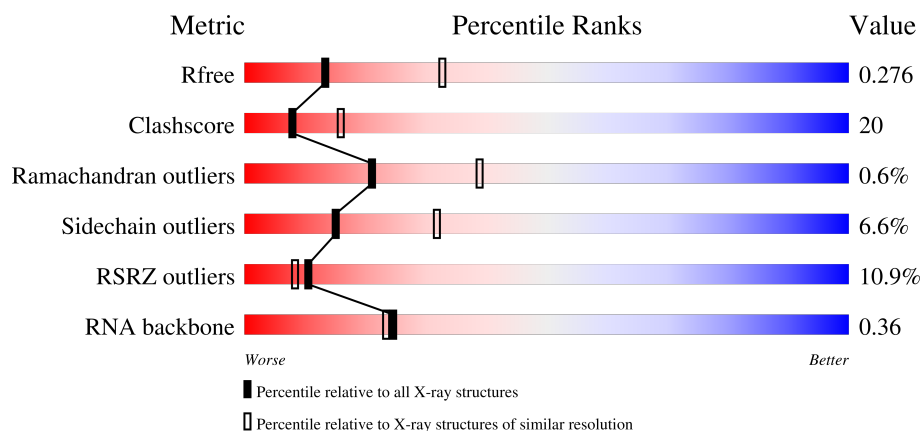
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.84 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}	180053	1520 (2.86-2.82)
Clashscore	190562	1559 (2.86-2.82)
Ramachandran outliers	187476	1517 (2.86-2.82)
Sidechain outliers	187428	1518 (2.86-2.82)
RSRZ outliers	180081	1521 (2.86-2.82)
RNA backbone	3983	1035 (3.06-2.62)

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 $\leq 5\%$. 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	9	33% 56% 33% 11%
1	D	9	33% 33% 56% 11%
2	A	335	5% 60% 35% ...
2	B	335	15% 58% 34% ...

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5653 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 RNA chain called HO endonuclease 3' UTR binding sequence.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	9	Total	C	N	O	P	0	0	0
			186	85	30	63	8			
1	D	9	Total	C	N	O	P	0	0	0
			186	85	30	63	8			

- Molecule 2 is a protein called Protein PUF4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	328	Total	C	N	O	S	0	0	0
			2611	1658	446	494	13			
2	B	323	Total	C	N	O	S	0	0	0
			2570	1631	439	487	13			

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0

- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	2	Total Na 2 2	0	0

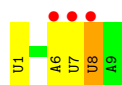
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	C	1	Total O 1 1	0	0
5	D	3	Total O 3 3	0	0
5	A	37	Total O 37 37	0	0
5	B	17	Total O 17 17	0	0

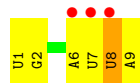
3 Residue-property plots [i](#)

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.

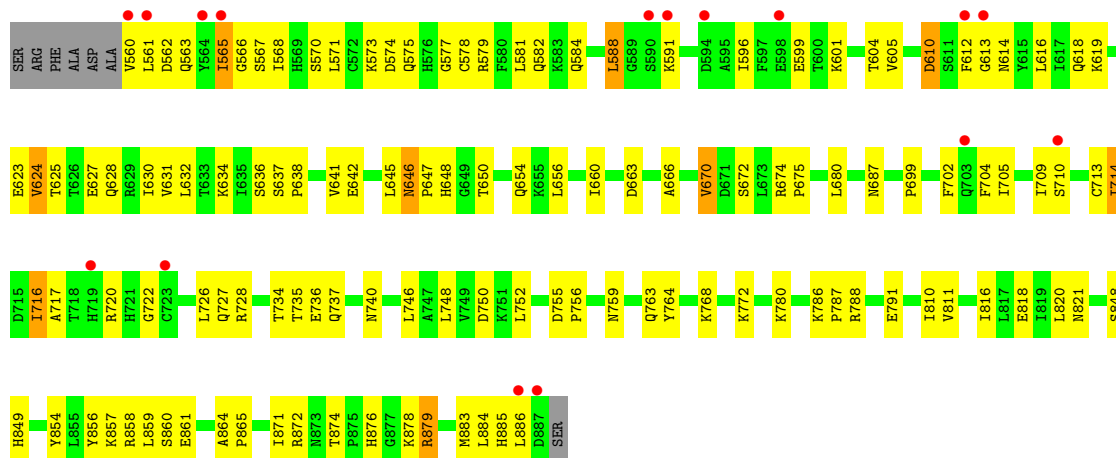
- Molecule 1: HO endonuclease 3' UTR binding sequence



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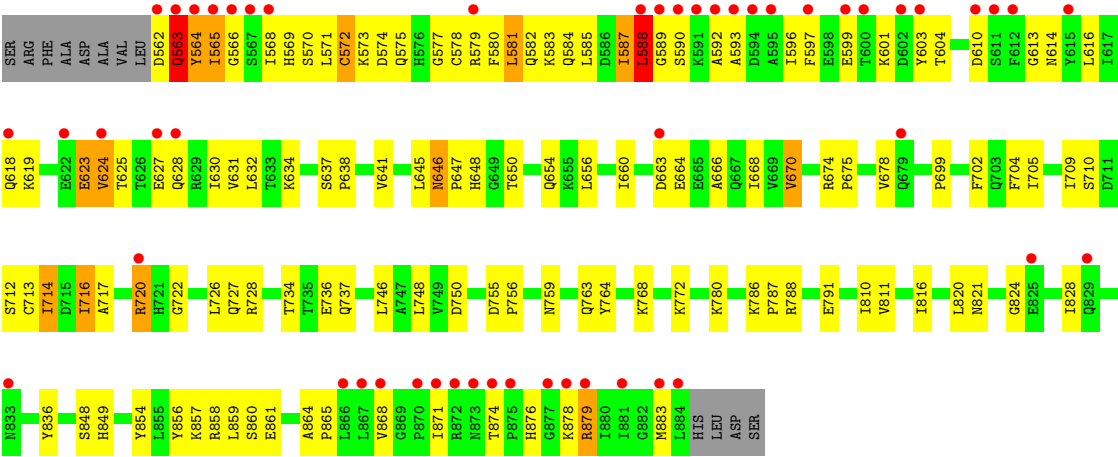


- Molecule 2: Protein PUF4



- Molecule 2: Protein PUF4





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	65.86Å 136.18Å 158.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	14.90 – 2.84 14.90 – 2.84	Depositor EDS
% Data completeness (in resolution range)	(Not available) (14.90-2.84) 92.8 (14.90-2.84)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.77 (at 2.86Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.240 , 0.280 0.239 , 0.276	Depositor DCC
R_{free} test set	1645 reflections (4.80%)	wwPDB-VP
Wilson B-factor (Å ²)	37.9	Xtriage
Anisotropy	0.242	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 61.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	5653	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.14% 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: NA, SO4

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	0.47	0/207	1.02	0/320
1	D	0.50	0/207	0.97	0/320
2	A	0.53	0/2655	0.88	3/3592 (0.1%)
2	B	0.50	0/2613	0.89	5/3534 (0.1%)
All	All	0.51	0/5682	0.89	8/7766 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	B	0	1

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	646	ASN	CA-C-N	7.50	127.38	119.05
2	B	646	ASN	C-N-CA	7.50	127.38	119.05
2	A	646	ASN	CA-C-N	7.38	127.24	119.05
2	A	646	ASN	C-N-CA	7.38	127.24	119.05
2	B	588	LEU	N-CA-C	-7.35	104.34	112.72
2	B	563	GLN	N-CA-C	5.92	116.37	108.38
2	A	588	LEU	N-CA-C	5.73	119.92	111.04
2	B	712	SER	N-CA-C	-5.69	103.75	111.55

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	588	LEU	Peptide

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	186	0	96	5	0
1	D	186	0	96	6	0
2	A	2611	0	2654	100	0
2	B	2570	0	2612	115	0
3	A	20	0	0	0	0
3	B	20	0	0	1	0
4	B	2	0	0	0	0
5	A	37	0	0	2	0
5	B	17	0	0	2	0
5	C	1	0	0	1	0
5	D	3	0	0	0	0
All	All	5653	0	5458	218	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 20.

All (218) 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:624:VAL:HG13	2:B:628:GLN:HB2	1.43	0.97
2:A:624:VAL:HG13	2:A:628:GLN:HB2	1.43	0.96
2:B:734:THR:HG22	2:B:736:GLU:H	1.29	0.96
2:A:734:THR:HG22	2:A:736:GLU:H	1.31	0.93
2:A:704:PHE:CZ	2:A:705:ILE:HG13	2.12	0.84
2:A:578:CYS:O	2:A:582:GLN:HG3	1.77	0.83
2:B:704:PHE:CZ	2:B:705:ILE:HG13	2.14	0.82
2:B:572:CYS:HB3	2:B:616:LEU:HD22	1.60	0.81
2:B:584:GLN:HB3	2:B:592:ALA:HB1	1.61	0.81
2:A:562:ASP:HB3	2:A:565:ILE:HG13	1.62	0.79
2:A:575:GLN:HB2	2:A:612:PHE:CE1	2.17	0.79
2:A:714:ILE:HD11	2:A:748:LEU:HD13	1.66	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:720:ARG:HG2	5:B:966:HOH:O	1.85	0.75
2:B:734:THR:HG22	2:B:736:GLU:N	2.01	0.75
2:B:570:SER:HA	2:B:573:LYS:HE2	1.67	0.74
2:B:714:ILE:HD11	2:B:748:LEU:HD13	1.69	0.74
2:B:604:THR:HG21	2:B:632:LEU:HD22	1.69	0.73
2:A:604:THR:HG21	2:A:632:LEU:HD22	1.70	0.73
2:B:568:ILE:HG23	2:B:599:GLU:HB3	1.71	0.72
2:B:624:VAL:CG1	2:B:628:GLN:HB2	2.20	0.71
2:A:810:ILE:HD12	2:A:811:VAL:HG13	1.71	0.71
2:A:734:THR:HG22	2:A:736:GLU:N	2.05	0.70
2:A:650:THR:O	2:A:654:GLN:HG3	1.91	0.70
1:D:1:U:N3	2:B:876:HIS:HB3	2.06	0.69
2:B:810:ILE:HD12	2:B:811:VAL:HG13	1.75	0.69
2:A:704:PHE:CE1	2:A:705:ILE:HG13	2.28	0.68
2:B:704:PHE:CE1	2:B:705:ILE:HG13	2.30	0.67
2:B:584:GLN:HB3	2:B:592:ALA:CB	2.25	0.66
2:A:575:GLN:HE21	2:A:579:ARG:HE	1.44	0.65
2:A:610:ASP:HB2	5:A:961:HOH:O	1.97	0.65
2:A:705:ILE:CG2	2:A:709:ILE:HD12	2.28	0.64
2:B:585:LEU:O	2:B:623:GLU:HG2	1.97	0.64
2:B:746:LEU:HD21	2:B:780:LYS:HB3	1.80	0.64
2:A:714:ILE:CD1	2:A:748:LEU:HD13	2.27	0.64
2:A:624:VAL:CG1	2:A:628:GLN:HB2	2.23	0.63
2:B:582:GLN:HG2	2:B:616:LEU:CD1	2.29	0.62
2:B:705:ILE:CG2	2:B:709:ILE:HD12	2.29	0.62
2:A:570:SER:HA	2:A:573:LYS:HE3	1.81	0.62
2:B:714:ILE:CD1	2:B:748:LEU:HD13	2.29	0.61
2:B:864:ALA:HB3	2:B:865:PRO:HD3	1.82	0.61
2:A:874:THR:O	2:A:878:LYS:HG3	2.00	0.61
2:B:650:THR:O	2:B:654:GLN:HG3	2.00	0.61
2:A:864:ALA:HB3	2:A:865:PRO:HD3	1.83	0.61
2:A:604:THR:HG21	2:A:632:LEU:CD2	2.30	0.61
2:B:582:GLN:HG2	2:B:616:LEU:HD13	1.83	0.61
2:B:816:ILE:HD13	2:B:848:SER:HB3	1.83	0.60
2:B:879:ARG:NH1	5:B:970:HOH:O	2.34	0.60
1:C:6:A:H4'	5:C:10:HOH:O	2.02	0.60
2:B:874:THR:O	2:B:878:LYS:HG3	2.02	0.60
2:B:759:ASN:O	2:B:763:GLN:HG3	2.01	0.60
2:B:705:ILE:HG22	2:B:709:ILE:HD12	1.83	0.59
2:B:604:THR:HG21	2:B:632:LEU:CD2	2.32	0.59
2:B:674:ARG:HB3	2:B:675:PRO:HD3	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:8:U:H5''	1:C:8:U:H6	1.68	0.59
2:B:575:GLN:NE2	2:B:579:ARG:HH21	2.01	0.59
2:A:562:ASP:CB	2:A:565:ILE:HG13	2.31	0.58
2:B:734:THR:HB	2:B:737:GLN:HG3	1.83	0.58
2:A:746:LEU:HD21	2:A:780:LYS:HB3	1.85	0.58
2:B:734:THR:HG23	3:B:953:SO4:O1	2.03	0.58
2:A:674:ARG:HB3	2:A:675:PRO:HD3	1.85	0.57
2:A:759:ASN:O	2:A:763:GLN:HG3	2.05	0.57
2:A:705:ILE:HG22	2:A:709:ILE:HD12	1.87	0.57
2:B:584:GLN:CB	2:B:592:ALA:HB1	2.34	0.56
2:A:567:SER:O	2:A:571:LEU:HD23	2.06	0.56
2:A:571:LEU:O	2:A:577:GLY:HA3	2.06	0.56
2:A:613:GLY:O	2:A:616:LEU:N	2.39	0.56
2:A:699:PRO:HA	2:A:702:PHE:CZ	2.40	0.56
1:D:8:U:H5''	1:D:8:U:H6	1.71	0.55
2:B:568:ILE:HG23	2:B:599:GLU:CB	2.35	0.55
2:B:820:LEU:HD21	2:B:859:LEU:HD23	1.89	0.55
2:A:716:ILE:HG23	2:A:722:GLY:CA	2.36	0.55
2:B:585:LEU:C	2:B:589:GLY:HA2	2.31	0.55
2:A:849:HIS:CE1	2:A:886:LEU:HB3	2.41	0.55
2:A:563:GLN:N	2:A:563:GLN:CD	2.66	0.54
2:A:638:PRO:HB3	2:B:678:VAL:HG11	1.90	0.54
2:A:857:LYS:O	2:A:861:GLU:HG2	2.08	0.54
2:B:572:CYS:HB3	2:B:616:LEU:CD2	2.33	0.54
2:B:587:ILE:HD12	2:B:588:LEU:H	1.73	0.54
2:B:714:ILE:CG1	2:B:748:LEU:HD13	2.37	0.54
2:A:601:LYS:HA	2:A:604:THR:HG23	1.90	0.54
2:A:656:LEU:C	2:A:656:LEU:HD23	2.33	0.53
2:B:699:PRO:HA	2:B:702:PHE:CZ	2.43	0.53
2:A:768:LYS:O	2:A:772:LYS:HG3	2.08	0.53
2:B:569:HIS:CD2	2:B:603:TYR:CZ	2.97	0.53
2:A:704:PHE:CZ	2:A:705:ILE:CG1	2.89	0.53
2:B:849:HIS:HB2	2:B:856:TYR:CD1	2.43	0.53
2:A:816:ILE:HD13	2:A:848:SER:HB3	1.91	0.53
2:A:641:VAL:O	2:A:645:LEU:HG	2.07	0.53
2:B:568:ILE:CG2	2:B:599:GLU:HB3	2.37	0.53
2:B:601:LYS:HA	2:B:604:THR:HG23	1.90	0.53
2:B:704:PHE:CZ	2:B:705:ILE:CG1	2.91	0.52
2:A:627:GLU:O	2:A:631:VAL:HG23	2.09	0.52
2:A:709:ILE:HG12	2:A:716:ILE:HD12	1.92	0.52
2:A:714:ILE:CG1	2:A:748:LEU:HD13	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:849:HIS:HB2	2:B:856:TYR:CE1	2.44	0.52
2:A:820:LEU:HD21	2:A:859:LEU:HD23	1.91	0.52
2:A:614:ASN:ND2	2:A:648:HIS:HB3	2.25	0.52
2:A:849:HIS:HB2	2:A:856:TYR:CD1	2.45	0.52
2:B:587:ILE:HD12	2:B:588:LEU:N	2.25	0.51
2:B:710:SER:O	2:B:713:CYS:HB2	2.09	0.51
2:A:710:SER:O	2:A:713:CYS:HB2	2.10	0.51
2:B:578:CYS:O	2:B:582:GLN:HG3	2.11	0.51
2:B:768:LYS:O	2:B:772:LYS:HG3	2.09	0.51
1:D:2:G:H1'	2:B:836:TYR:CD1	2.46	0.51
2:A:699:PRO:HA	2:A:702:PHE:CE2	2.46	0.51
2:B:755:ASP:OD1	2:B:756:PRO:HD2	2.11	0.51
2:A:849:HIS:HB2	2:A:856:TYR:CE1	2.46	0.50
2:B:569:HIS:CE1	2:B:603:TYR:CE1	3.00	0.50
2:A:666:ALA:O	2:A:670:VAL:HG13	2.11	0.50
2:A:734:THR:HB	2:A:737:GLN:HG3	1.92	0.50
2:A:642:GLU:HB3	2:B:714:ILE:HB	1.94	0.50
2:B:641:VAL:O	2:B:645:LEU:HG	2.12	0.50
2:B:613:GLY:O	2:B:616:LEU:N	2.45	0.50
2:A:601:LYS:O	2:A:604:THR:HG23	2.12	0.49
2:B:716:ILE:HG23	2:B:722:GLY:CA	2.42	0.49
2:B:709:ILE:HG12	2:B:716:ILE:HD12	1.93	0.49
2:A:561:LEU:N	2:A:561:LEU:HD12	2.27	0.49
2:B:584:GLN:O	2:B:589:GLY:HA2	2.12	0.49
2:B:614:ASN:ND2	2:B:648:HIS:HB3	2.27	0.49
2:B:656:LEU:O	2:B:660:ILE:HG23	2.13	0.49
2:A:656:LEU:O	2:A:660:ILE:HG23	2.13	0.49
2:B:857:LYS:O	2:B:861:GLU:HG2	2.13	0.49
2:B:568:ILE:HD13	2:B:599:GLU:CD	2.38	0.48
2:B:624:VAL:CG1	2:B:625:THR:N	2.76	0.48
2:B:568:ILE:HG23	2:B:599:GLU:CG	2.44	0.48
2:A:614:ASN:O	2:A:618:GLN:HG3	2.13	0.48
2:A:755:ASP:OD1	2:A:756:PRO:HD2	2.13	0.48
2:B:578:CYS:SG	2:B:616:LEU:HB2	2.53	0.48
2:B:656:LEU:C	2:B:656:LEU:HD23	2.38	0.48
2:A:885:HIS:C	2:A:886:LEU:HD12	2.39	0.48
1:D:2:G:H1'	2:B:836:TYR:HD1	1.77	0.47
2:A:637:SER:HB2	2:A:672:SER:OG	2.14	0.47
1:C:6:A:C8	1:C:6:A:H3'	2.49	0.47
2:A:568:ILE:HB	2:A:599:GLU:OE1	2.14	0.47
2:B:727:GLN:HG2	2:B:764:TYR:CD2	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:6:A:C8	1:D:6:A:H3'	2.49	0.47
2:A:704:PHE:CG	2:A:705:ILE:N	2.83	0.47
2:A:646:ASN:HA	2:A:647:PRO:HD3	1.77	0.47
2:A:734:THR:HG22	2:A:735:THR:N	2.29	0.47
2:B:666:ALA:O	2:B:670:VAL:HG13	2.14	0.47
2:A:705:ILE:HG23	2:A:709:ILE:HD12	1.98	0.46
2:B:704:PHE:CG	2:B:705:ILE:N	2.83	0.46
2:B:786:LYS:HB3	2:B:787:PRO:HD3	1.97	0.46
2:A:614:ASN:HD22	2:A:648:HIS:HB3	1.80	0.46
2:A:630:ILE:O	2:A:634:LYS:HG3	2.16	0.46
2:B:585:LEU:CD2	2:B:593:ALA:HA	2.45	0.46
2:B:627:GLU:O	2:B:631:VAL:HG23	2.15	0.46
2:B:788:ARG:HD2	2:B:791:GLU:CD	2.39	0.46
2:B:637:SER:OG	2:B:638:PRO:HD3	2.15	0.46
2:A:565:ILE:CG2	2:A:566:GLY:N	2.78	0.46
2:B:563:GLN:O	2:B:564:TYR:CD1	2.69	0.46
2:B:585:LEU:O	2:B:589:GLY:CA	2.64	0.46
2:B:585:LEU:HG	2:B:596:ILE:HD12	1.96	0.46
2:A:788:ARG:HD2	2:A:791:GLU:CD	2.42	0.45
2:B:714:ILE:HG12	2:B:748:LEU:HD13	1.97	0.45
2:A:568:ILE:HD13	2:A:596:ILE:HG12	1.99	0.45
2:B:574:ASP:O	2:B:578:CYS:HB2	2.15	0.45
2:A:637:SER:OG	2:A:638:PRO:HD3	2.16	0.45
2:B:585:LEU:O	2:B:623:GLU:CG	2.64	0.45
2:B:624:VAL:HG13	2:B:625:THR:N	2.31	0.45
2:B:646:ASN:HA	2:B:647:PRO:HD3	1.76	0.45
2:B:564:TYR:C	2:B:564:TYR:CD2	2.95	0.45
2:B:824:GLY:O	2:B:828:ILE:HG13	2.16	0.45
1:C:8:U:O2	2:A:614:ASN:ND2	2.50	0.44
2:A:714:ILE:HG12	2:A:748:LEU:HD13	1.99	0.44
2:B:630:ILE:O	2:B:634:LYS:HG3	2.17	0.44
2:B:699:PRO:HA	2:B:702:PHE:CE2	2.51	0.44
2:B:727:GLN:HG2	2:B:764:TYR:CG	2.53	0.44
2:B:577:GLY:O	2:B:581:LEU:HB2	2.17	0.44
2:A:734:THR:CG2	2:A:735:THR:N	2.80	0.43
2:A:748:LEU:O	2:A:752:LEU:HG	2.18	0.43
2:B:582:GLN:HG2	2:B:616:LEU:HD12	2.00	0.43
2:A:849:HIS:NE2	2:A:886:LEU:HB3	2.33	0.43
2:B:717:ALA:HB2	2:B:726:LEU:CD1	2.48	0.43
2:A:578:CYS:SG	2:A:612:PHE:HB2	2.58	0.43
2:B:565:ILE:HG22	2:B:566:GLY:N	2.32	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:624:VAL:CG1	2:A:625:THR:N	2.81	0.43
2:A:854:TYR:O	2:A:858:ARG:HG2	2.19	0.43
2:A:879:ARG:O	2:A:883:MET:HG3	2.19	0.43
2:A:728:ARG:HD3	2:A:728:ARG:HA	1.85	0.43
2:B:601:LYS:O	2:B:604:THR:HG23	2.18	0.43
2:A:786:LYS:HB3	2:A:787:PRO:HD3	1.99	0.43
2:B:854:TYR:O	2:B:858:ARG:HG2	2.19	0.43
2:A:591:LYS:NZ	5:A:987:HOH:O	2.45	0.43
2:B:585:LEU:O	2:B:589:GLY:HA3	2.19	0.43
2:B:664:GLU:O	2:B:668:ILE:HG13	2.19	0.43
2:A:858:ARG:HA	2:A:858:ARG:HD3	1.79	0.43
2:B:562:ASP:CG	2:B:563:GLN:H	2.27	0.43
2:B:585:LEU:HB3	2:B:623:GLU:HG2	2.00	0.43
2:B:614:ASN:O	2:B:618:GLN:HG3	2.18	0.42
1:D:1:U:C2	2:B:876:HIS:HB3	2.54	0.42
2:A:884:LEU:HD23	2:A:884:LEU:HA	1.82	0.42
2:B:619:LYS:HA	2:B:619:LYS:HD2	1.91	0.42
2:A:818:GLU:OE1	2:A:818:GLU:HA	2.20	0.42
2:B:585:LEU:HD21	2:B:596:ILE:HD12	2.01	0.42
2:A:619:LYS:HD2	2:A:619:LYS:HA	1.87	0.42
2:A:727:GLN:HG2	2:A:764:TYR:CD2	2.54	0.42
2:A:612:PHE:O	2:A:613:GLY:C	2.63	0.42
2:A:581:LEU:O	2:A:584:GLN:HB2	2.20	0.42
2:B:580:PHE:O	2:B:583:LYS:HB2	2.20	0.42
2:B:728:ARG:HD3	2:B:728:ARG:HA	1.85	0.42
2:A:624:VAL:HG13	2:A:625:THR:N	2.36	0.41
2:A:849:HIS:HE1	2:A:886:LEU:O	2.04	0.41
2:B:858:ARG:HA	2:B:858:ARG:HD3	1.81	0.41
2:B:717:ALA:HA	2:B:726:LEU:HD12	2.03	0.41
2:A:717:ALA:HA	2:A:726:LEU:HD12	2.03	0.41
2:A:849:HIS:CE1	2:A:886:LEU:O	2.74	0.41
2:B:585:LEU:CD2	2:B:596:ILE:HD12	2.51	0.41
2:B:585:LEU:CG	2:B:596:ILE:HD12	2.51	0.41
2:B:588:LEU:O	2:B:590:SER:N	2.54	0.41
2:A:872:ARG:HH11	2:A:872:ARG:HD2	1.75	0.41
2:B:581:LEU:O	2:B:596:ILE:HD11	2.21	0.40
2:A:588:LEU:HD23	2:A:588:LEU:HA	1.71	0.40
2:A:849:HIS:CE1	2:A:886:LEU:CB	3.04	0.40
2:A:883:MET:HA	2:A:886:LEU:HD11	2.03	0.40
2:A:768:LYS:HB3	2:A:768:LYS:HE2	1.85	0.40
2:B:879:ARG:O	2:B:883:MET:HG3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1:U:N3	2:A:876:HIS:HB3	2.36	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	
2	A	326/335 (97%)	309 (95%)	15 (5%)	2 (1%)	21	39
2	B	321/335 (96%)	301 (94%)	18 (6%)	2 (1%)	21	39
All	All	647/670 (97%)	610 (94%)	33 (5%)	4 (1%)	21	39

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	821	ASN
2	A	565	ILE
2	A	821	ASN
2	B	597	PHE

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	
2	A	298/303 (98%)	279 (94%)	19 (6%)	16	33

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	293/303 (97%)	273 (93%)	20 (7%)	14	31
All	All	591/606 (98%)	552 (93%)	39 (7%)	15	32

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	560	VAL
2	A	574	ASP
2	A	605	VAL
2	A	610	ASP
2	A	623	GLU
2	A	624	VAL
2	A	636	SER
2	A	663	ASP
2	A	670	VAL
2	A	680	LEU
2	A	687	ASN
2	A	714	ILE
2	A	716	ILE
2	A	720	ARG
2	A	740	ASN
2	A	750	ASP
2	A	860	SER
2	A	871	ILE
2	A	879	ARG
2	B	563	GLN
2	B	564	TYR
2	B	565	ILE
2	B	571	LEU
2	B	572	CYS
2	B	581	LEU
2	B	587	ILE
2	B	610	ASP
2	B	623	GLU
2	B	624	VAL
2	B	663	ASP
2	B	670	VAL
2	B	714	ILE
2	B	716	ILE
2	B	720	ARG
2	B	750	ASP
2	B	860	SER

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Mol	Chain	Res	Type
2	B	868	VAL
2	B	871	ILE
2	B	879	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	575	GLN
2	A	740	ASN
2	A	773	ASN
2	A	800	ASN
2	A	849	HIS
2	A	885	HIS
2	B	569	HIS
2	B	575	GLN
2	B	639	HIS
2	B	688	HIS
2	B	740	ASN
2	B	800	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	C	8/9 (88%)	2 (25%)	0
1	D	8/9 (88%)	3 (37%)	0
All	All	16/18 (88%)	5 (31%)	0

All (5) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	C	7	U
1	C	8	U
1	D	7	U
1	D	8	U
1	D	9	A

There are no RNA pucker outliers to report.

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 2 are monoatomic - leaving 8 for Mogul analysis.

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	$\# Z > 2$
3	SO4	B	952	-	4,4,4	0.24	0	6,6,6	0.13	0
3	SO4	A	956	-	4,4,4	0.20	0	6,6,6	0.13	0
3	SO4	A	951	-	4,4,4	0.27	0	6,6,6	0.23	0
3	SO4	B	955	-	4,4,4	0.23	0	6,6,6	0.09	0
3	SO4	A	958	-	4,4,4	0.62	0	6,6,6	0.28	0
3	SO4	A	957	-	4,4,4	0.24	0	6,6,6	0.15	0
3	SO4	B	953	-	4,4,4	0.22	0	6,6,6	0.15	0
3	SO4	B	954	-	4,4,4	0.27	0	6,6,6	0.12	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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	953	SO4	1	0

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

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	9/9 (100%)	1.40	3 (33%) 1 0	46, 55, 104, 157	0
1	D	9/9 (100%)	1.26	3 (33%) 1 0	48, 66, 94, 159	0
2	A	328/335 (97%)	0.06	16 (4%) 35 28	12, 28, 69, 111	0
2	B	323/335 (96%)	0.61	51 (15%) 5 3	13, 35, 93, 133	0
All	All	669/688 (97%)	0.36	73 (10%) 10 8	12, 32, 88, 159	0

All (73) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	592	ALA	9.5
2	B	590	SER	6.5
2	B	881	ILE	6.2
2	A	560	VAL	5.0
2	B	599	GLU	5.0
2	B	868	VAL	4.7
2	B	603	TYR	4.7
2	B	567	SER	4.6
2	A	561	LEU	4.5
2	B	564	TYR	4.5
2	B	874	THR	4.4
2	B	589	GLY	4.3
2	B	563	GLN	4.3
2	B	612	PHE	4.2
2	B	877	GLY	4.2
2	B	593	ALA	4.1
1	C	7	U	4.0
2	B	562	ASP	4.0
2	B	873	ASN	4.0
2	B	872	ARG	3.9
2	B	568	ILE	3.9

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Mol	Chain	Res	Type	RSRZ
1	C	8	U	3.8
2	B	600	THR	3.7
2	B	627	GLU	3.6
2	B	591	LYS	3.6
2	B	866	LEU	3.6
2	A	887	ASP	3.6
2	B	602	ASP	3.6
2	A	590	SER	3.5
2	B	610	ASP	3.5
2	A	591	LYS	3.4
2	B	566	GLY	3.3
2	A	613	GLY	3.3
2	B	825	GLU	3.3
2	A	703	GLN	3.2
2	B	622	GLU	3.2
2	B	618	GLN	3.1
2	B	871	ILE	3.1
1	D	7	U	3.0
2	B	595	ALA	3.0
2	B	883	MET	3.0
2	B	875	PRO	2.9
2	B	829	GLN	2.9
1	D	6	A	2.9
2	B	870	PRO	2.8
2	B	594	ASP	2.8
2	A	710	SER	2.8
1	D	8	U	2.8
2	A	612	PHE	2.8
2	B	679	GLN	2.8
2	B	867	LEU	2.7
2	B	611	SER	2.7
1	C	6	A	2.7
2	A	564	TYR	2.6
2	B	884	LEU	2.6
2	A	598	GLU	2.6
2	B	628	GLN	2.6
2	B	579	ARG	2.5
2	B	663	ASP	2.5
2	A	723	CYS	2.5
2	B	833	ASN	2.4
2	B	720	ARG	2.3
2	A	565	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
2	B	565	ILE	2.3
2	B	597	PHE	2.2
2	A	886	LEU	2.1
2	B	615	TYR	2.1
2	A	719	HIS	2.1
2	B	624	VAL	2.1
2	A	594	ASP	2.1
2	B	878	LYS	2.1
2	B	588	LEU	2.1
2	B	879	ARG	2.1

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 oligosaccharides 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(Å ²)	Q<0.9
3	SO4	B	952	5/5	0.59	0.18	134,135,137,142	0
3	SO4	B	954	5/5	0.75	0.21	100,101,111,128	0
3	SO4	A	957	5/5	0.82	0.17	113,113,124,127	0
3	SO4	B	953	5/5	0.83	0.14	104,105,109,120	0
4	NA	B	901	1/1	0.87	0.21	30,30,30,30	0
3	SO4	A	956	5/5	0.93	0.15	95,106,113,117	0
3	SO4	B	955	5/5	0.94	0.14	109,112,114,118	0
3	SO4	A	958	5/5	0.95	0.28	30,30,30,30	0
3	SO4	A	951	5/5	0.96	0.08	41,55,72,73	0
4	NA	B	902	1/1	0.97	0.23	30,30,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.