



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:35 AM UTC

PDB ID : 3BX3 / pdb_00003bx3
Title : Puf4 T650C/C724R Mutant bound to Cox17 RNA 3' UTR recognition sequence
Authors : Miller, M.T.; Higgin, J.J.; Hall, T.M.T.
Deposited on : 2008-01-11
Resolution : 3.00 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

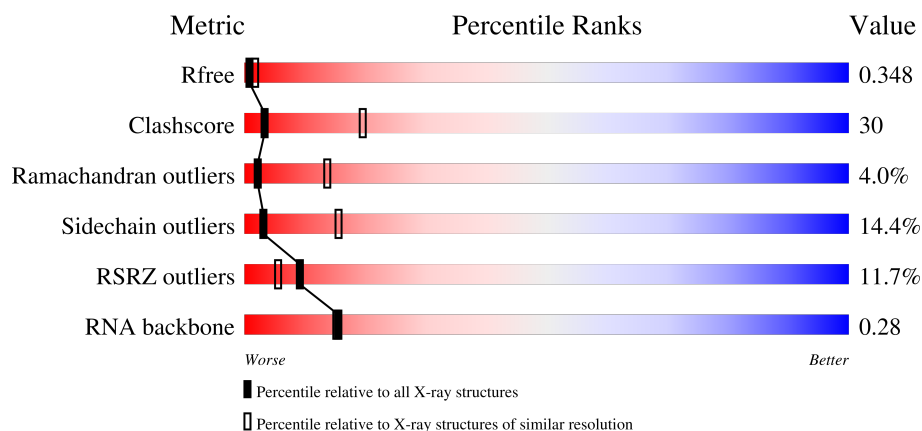
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)
RNA backbone	3983	1109 (3.20-2.80)

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	8	62% 50% 38% 12%
1	D	8	50% 38% 12%
2	A	335	12% 38% 45% 13% ••
2	B	335	8% 42% 37% 10% • 10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5349 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 COX17 RNA target sequence.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	8	Total	C	N	O	P	0	0	0
			166	76	28	55	7			
1	D	8	Total	C	N	O	P	0	0	0
			166	76	28	55	7			

- Molecule 2 is a protein called Protein PUF4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	325	Total	C	N	O	S	0	0	0
			2592	1645	446	488	13			
2	B	300	Total	C	N	O	S	0	0	0
			2380	1511	408	450	11			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	650	CYS	THR	engineered mutation	UNP P25339
A	724	ARG	CYS	engineered mutation	UNP P25339
B	650	CYS	THR	engineered mutation	UNP P25339
B	724	ARG	CYS	engineered mutation	UNP P25339

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

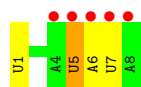


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

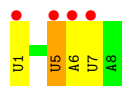
3 Residue-property plots

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.

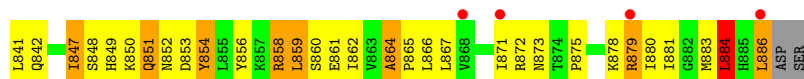
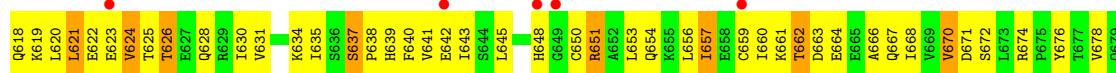
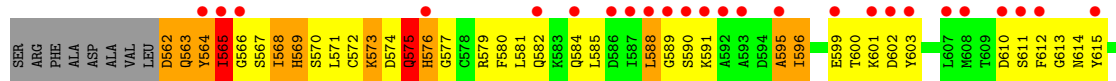
- Molecule 1: COX17 RNA target sequence



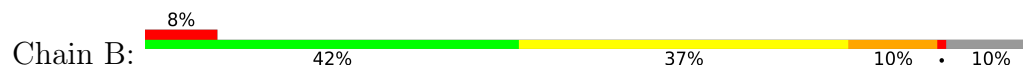
- Molecule 1: COX17 RNA target sequence

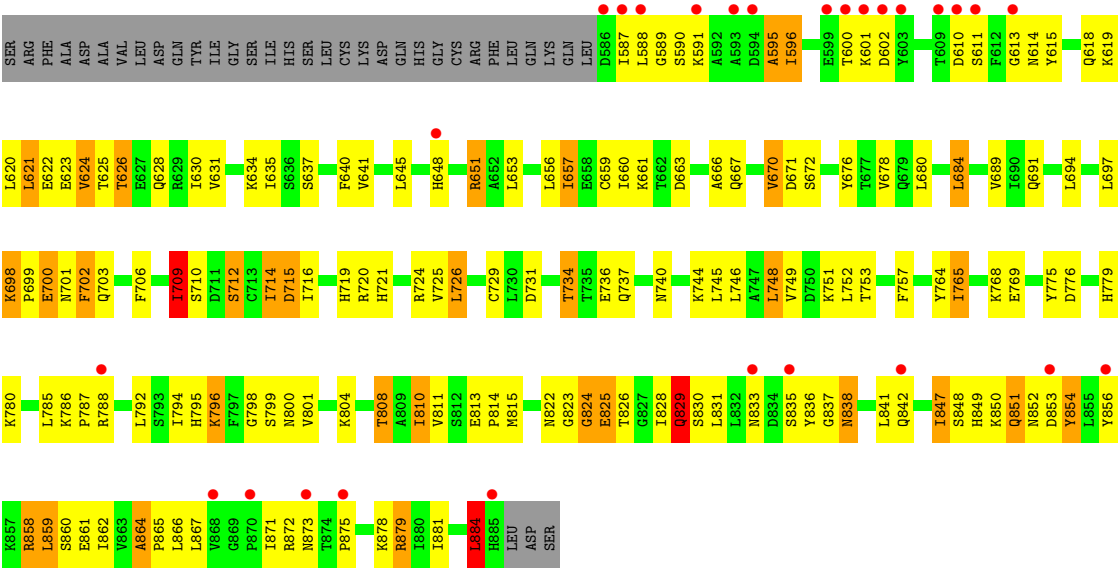


- Molecule 2: Protein PUF4



- 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	66.06Å 137.12Å 160.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.40 – 3.00 24.40 – 3.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (24.40-3.00) 98.0 (24.40-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.14 (at 2.99Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.290 , 0.350 0.292 , 0.348	Depositor DCC
R_{free} test set	1500 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	38.1	Xtriage
Anisotropy	1.027	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 57.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.82	EDS
Total number of atoms	5349	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 22.41 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.7354e-03.*

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: 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	0.58	0/185	1.23	2/286 (0.7%)
1	D	0.61	0/185	1.20	2/286 (0.7%)
2	A	0.65	0/2636	1.10	21/3564 (0.6%)
2	B	0.61	0/2420	1.03	12/3275 (0.4%)
All	All	0.63	0/5426	1.08	37/7411 (0.5%)

There are no bond length outliers.

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	575	GLN	N-CA-C	-8.59	101.94	114.39
2	A	825	GLU	N-CA-C	-8.53	102.65	113.23
2	B	651	ARG	NE-CZ-NH2	7.59	126.03	119.20
2	B	651	ARG	NE-CZ-NH1	-7.08	114.42	121.50
2	A	698	LYS	CA-C-N	6.91	127.48	119.47
2	A	698	LYS	C-N-CA	6.91	127.48	119.47
2	B	884	LEU	N-CA-C	6.91	118.81	111.28
2	B	698	LYS	CA-C-N	6.61	127.13	119.47
2	B	698	LYS	C-N-CA	6.61	127.13	119.47
2	B	825	GLU	N-CA-C	-6.61	105.04	113.23
2	A	884	LEU	N-CA-C	6.51	118.38	111.28
2	A	864	ALA	CA-C-N	6.25	126.78	120.04
2	A	864	ALA	C-N-CA	6.25	126.78	120.04
2	A	651	ARG	NE-CZ-NH2	-6.18	113.63	119.20
2	A	577	GLY	N-CA-C	-6.15	107.27	115.32
2	A	651	ARG	CD-NE-CZ	5.91	132.67	124.40
2	A	637	SER	CA-C-N	-5.81	114.90	120.83
2	A	637	SER	C-N-CA	-5.81	114.90	120.83
2	A	564	TYR	N-CA-C	5.77	115.12	108.49
1	D	5	U	C1'-O4'-C4'	-5.69	104.21	109.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	864	ALA	CA-C-N	5.63	125.72	119.87
2	B	864	ALA	C-N-CA	5.63	125.72	119.87
2	A	714	ILE	N-CA-C	5.46	116.19	110.62
2	B	651	ARG	CD-NE-CZ	5.45	132.03	124.40
2	A	674	ARG	CA-C-N	-5.38	114.55	119.82
2	A	674	ARG	C-N-CA	-5.38	114.55	119.82
1	C	5	U	C1'-O4'-C4'	-5.37	104.53	109.90
2	A	880	ILE	CB-CA-C	-5.29	105.10	112.04
2	A	789	ALA	N-CA-C	5.23	116.67	111.07
2	A	668	ILE	N-CA-C	-5.18	106.75	111.67
2	A	709	ILE	N-CA-C	5.17	115.28	110.42
2	A	565	ILE	N-CA-C	5.15	120.05	109.34
1	C	5	U	C4'-C3'-C2'	-5.13	97.47	102.60
2	B	709	ILE	CB-CA-C	-5.11	105.28	111.87
1	D	5	U	C4'-C3'-C2'	-5.08	97.52	102.60
2	B	709	ILE	N-CA-C	5.01	115.13	110.42
2	B	829	GLN	N-CA-C	-5.00	105.73	111.14

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	166	0	86	11	0
1	D	166	0	86	8	0
2	A	2592	0	2636	196	0
2	B	2380	0	2415	142	0
3	A	25	0	0	1	0
3	B	20	0	0	1	0
All	All	5349	0	5223	321	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 30.

All (321) 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:A:642:GLU:HG3	2:B:714:ILE:CG2	1.60	1.31
2:B:753:THR:HG23	2:B:801:VAL:HG21	1.31	1.12
2:A:642:GLU:HG3	2:B:714:ILE:HG22	1.09	1.09
2:A:753:THR:HG23	2:A:801:VAL:HG21	1.31	1.07
2:B:734:THR:HG22	2:B:736:GLU:H	1.21	1.04
2:A:734:THR:HG22	2:A:736:GLU:H	1.20	1.02
2:B:808:THR:HG22	2:B:811:VAL:HG22	1.47	0.97
2:B:714:ILE:HD13	2:B:748:LEU:HD13	1.47	0.97
2:A:642:GLU:CG	2:B:714:ILE:HG22	1.94	0.97
2:A:714:ILE:HD13	2:A:748:LEU:HD13	1.51	0.93
2:A:808:THR:HG22	2:A:811:VAL:HG22	1.48	0.92
1:C:1:U:H3	2:A:838:ASN:ND2	1.67	0.92
2:A:575:GLN:HE22	2:A:579:ARG:HE	1.12	0.90
1:C:5:U:N3	2:A:724:ARG:HD3	1.89	0.85
2:A:572:CYS:O	2:A:573:LYS:HG3	1.76	0.85
2:A:734:THR:HG22	2:A:736:GLU:N	1.91	0.85
1:C:1:U:H3	2:A:838:ASN:HD21	1.25	0.85
2:A:753:THR:HG23	2:A:801:VAL:CG2	2.09	0.83
2:B:734:THR:HG22	2:B:736:GLU:N	1.93	0.82
2:A:588:LEU:HB3	2:A:589:GLY:CA	2.09	0.82
2:B:753:THR:HG23	2:B:801:VAL:CG2	2.09	0.81
2:A:849:HIS:HE1	2:A:886:LEU:C	1.89	0.80
2:A:883:MET:O	2:A:886:LEU:HD22	1.85	0.77
1:D:1:U:H3	2:B:838:ASN:HD21	1.33	0.77
2:B:864:ALA:HB3	2:B:865:PRO:HD3	1.65	0.76
2:A:621:LEU:N	2:A:621:LEU:HD23	2.01	0.74
2:A:676:TYR:OH	2:B:719:HIS:HD2	1.69	0.74
2:A:642:GLU:HG3	2:B:714:ILE:HG21	1.63	0.74
2:A:735:THR:HG21	3:A:905:SO4:O2	1.88	0.74
2:B:621:LEU:HD23	2:B:621:LEU:N	2.03	0.74
2:A:864:ALA:HB3	2:A:865:PRO:HD3	1.68	0.74
1:D:5:U:N3	2:B:724:ARG:HD3	2.03	0.73
2:A:582:GLN:HE22	2:A:615:TYR:HB3	1.54	0.73
2:A:588:LEU:HB3	2:A:589:GLY:HA2	1.71	0.72
2:A:794:ILE:O	2:A:836:TYR:HD2	1.74	0.71
2:B:808:THR:HG22	2:B:811:VAL:CG2	2.20	0.71
2:A:588:LEU:CB	2:A:589:GLY:HA2	2.20	0.70
2:A:588:LEU:HD22	2:A:589:GLY:HA2	1.73	0.70
1:C:7:U:H3	2:A:618:GLN:HE22	1.40	0.69
2:B:794:ILE:O	2:B:836:TYR:HD2	1.76	0.69
2:A:849:HIS:CE1	2:A:886:LEU:HB3	2.28	0.69
2:B:734:THR:CG2	2:B:736:GLU:H	2.01	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:808:THR:HG22	2:A:811:VAL:CG2	2.21	0.68
2:A:667:GLN:O	2:A:671:ASP:OD1	2.13	0.67
2:A:849:HIS:CE1	2:A:886:LEU:CB	2.78	0.67
2:A:571:LEU:HD23	2:A:571:LEU:N	2.09	0.67
2:B:678:VAL:HG22	2:B:716:ILE:HD11	1.74	0.67
2:A:678:VAL:HG22	2:A:716:ILE:HD11	1.77	0.66
1:C:5:U:C2	2:A:724:ARG:HD3	2.30	0.65
2:A:715:ASP:O	2:A:719:HIS:HB2	1.98	0.64
2:B:611:SER:HA	2:B:648:HIS:CE1	2.32	0.64
2:A:588:LEU:CD2	2:A:589:GLY:HA2	2.28	0.63
2:B:813:GLU:HB2	2:B:814:PRO:HD3	1.81	0.63
2:A:700:GLU:CD	2:A:700:GLU:H	2.07	0.63
2:B:610:ASP:OD2	2:B:613:GLY:N	2.32	0.63
2:B:829:GLN:HG2	2:B:866:LEU:HD13	1.80	0.63
2:B:700:GLU:CD	2:B:700:GLU:H	2.05	0.63
2:A:611:SER:HA	2:A:648:HIS:CE1	2.34	0.62
2:B:715:ASP:OD1	2:B:715:ASP:N	2.32	0.62
2:A:564:TYR:CG	2:A:565:ILE:N	2.66	0.62
2:B:831:LEU:HD23	2:B:841:LEU:HB2	1.80	0.62
2:A:734:THR:CG2	2:A:736:GLU:H	2.03	0.62
2:B:734:THR:H	2:B:737:GLN:HB2	1.64	0.62
2:A:574:ASP:HB3	2:A:576:HIS:HD2	1.66	0.61
2:B:715:ASP:O	2:B:719:HIS:HB2	2.01	0.61
2:B:656:LEU:C	2:B:656:LEU:HD23	2.26	0.61
2:A:610:ASP:OD2	2:A:613:GLY:N	2.34	0.60
2:A:676:TYR:OH	2:B:719:HIS:CD2	2.52	0.60
2:A:625:THR:HG23	2:A:628:GLN:OE1	2.02	0.60
2:B:776:ASP:O	2:B:780:LYS:HG3	2.01	0.59
2:B:875:PRO:O	2:B:879:ARG:HG2	2.01	0.59
2:A:656:LEU:C	2:A:656:LEU:HD23	2.27	0.59
2:A:829:GLN:HE22	2:A:833:ASN:ND2	2.00	0.59
2:B:667:GLN:O	2:B:671:ASP:OD1	2.21	0.58
2:A:829:GLN:HG2	2:A:866:LEU:HD13	1.84	0.58
2:A:715:ASP:OD1	2:A:715:ASP:N	2.33	0.58
2:A:813:GLU:HB2	2:A:814:PRO:HD3	1.85	0.58
2:B:698:LYS:HE3	2:B:700:GLU:OE1	2.03	0.58
1:D:1:U:H3	2:B:838:ASN:ND2	1.99	0.58
2:A:699:PRO:HA	2:A:702:PHE:CE2	2.39	0.58
2:A:776:ASP:O	2:A:780:LYS:HG3	2.03	0.58
2:A:875:PRO:O	2:A:879:ARG:HG2	2.03	0.58
2:A:734:THR:H	2:A:737:GLN:HB2	1.68	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:822:ASN:C	2:A:824:GLY:H	2.12	0.57
2:B:823:GLY:O	2:B:824:GLY:O	2.22	0.57
2:B:625:THR:HG23	2:B:628:GLN:OE1	2.03	0.57
2:A:640:PHE:HD2	2:A:672:SER:HB3	1.68	0.57
2:B:698:LYS:O	2:B:701:ASN:HB2	2.04	0.57
2:B:829:GLN:HE22	2:B:833:ASN:ND2	2.02	0.57
1:C:7:U:C4	2:A:651:ARG:HD2	2.40	0.56
2:A:584:GLN:O	2:A:588:LEU:HB2	2.05	0.56
2:A:614:ASN:O	2:A:618:GLN:HG3	2.06	0.56
2:A:788:ARG:NH2	2:A:792:LEU:HD21	2.21	0.56
2:A:831:LEU:HD23	2:A:841:LEU:HB2	1.86	0.56
2:A:698:LYS:O	2:A:701:ASN:HB2	2.05	0.55
2:B:858:ARG:NH1	2:B:861:GLU:OE2	2.39	0.55
2:A:719:HIS:O	2:A:720:ARG:C	2.50	0.55
2:A:779:HIS:NE2	2:A:810:ILE:HD12	2.22	0.55
2:B:699:PRO:HA	2:B:702:PHE:CE2	2.41	0.55
2:A:574:ASP:HB3	2:A:576:HIS:CD2	2.42	0.55
2:A:849:HIS:CE1	2:A:886:LEU:C	2.79	0.55
2:A:858:ARG:NH1	2:A:861:GLU:OE2	2.40	0.55
2:B:749:VAL:HA	2:B:752:LEU:HB2	1.88	0.55
2:B:796:LYS:HE2	2:B:836:TYR:CE2	2.41	0.55
2:B:822:ASN:C	2:B:824:GLY:H	2.14	0.55
2:A:626:THR:O	2:A:630:ILE:HG13	2.07	0.54
2:A:564:TYR:CE2	2:A:565:ILE:HB	2.42	0.54
2:A:588:LEU:CB	2:A:589:GLY:CA	2.80	0.54
2:B:640:PHE:HD2	2:B:672:SER:HB3	1.72	0.54
1:D:7:U:C4	2:B:651:ARG:HD3	2.43	0.54
2:B:788:ARG:NH2	2:B:792:LEU:HD21	2.22	0.54
1:C:5:U:C2	2:A:724:ARG:CD	2.91	0.54
2:B:858:ARG:HH11	2:B:858:ARG:HG3	1.71	0.54
2:A:645:LEU:HD13	2:B:751:LYS:HD2	1.90	0.54
2:A:619:LYS:HA	2:A:622:GLU:HG3	1.90	0.53
2:A:676:TYR:HB2	2:A:680:LEU:HD12	1.90	0.53
2:A:698:LYS:HE3	2:A:700:GLU:OE1	2.08	0.53
2:A:749:VAL:HA	2:A:752:LEU:HB2	1.89	0.53
2:B:776:ASP:OD1	2:B:780:LYS:HE3	2.08	0.53
2:A:568:ILE:O	2:A:569:HIS:C	2.52	0.53
2:B:858:ARG:NH1	2:B:858:ARG:HG3	2.24	0.53
2:B:619:LYS:HA	2:B:622:GLU:HG3	1.91	0.52
2:A:706:PHE:CZ	2:A:729:CYS:HB3	2.45	0.52
2:A:719:HIS:O	2:A:721:HIS:N	2.42	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:564:TYR:CZ	2:A:565:ILE:HB	2.44	0.52
2:A:590:SER:C	2:A:591:LYS:HD3	2.35	0.52
2:A:666:ALA:O	2:A:670:VAL:HG13	2.10	0.52
2:A:776:ASP:OD1	2:A:780:LYS:HE3	2.10	0.52
2:A:796:LYS:HE2	2:A:836:TYR:CE2	2.44	0.52
2:B:626:THR:O	2:B:630:ILE:HG13	2.10	0.52
2:A:676:TYR:O	2:A:680:LEU:HD12	2.10	0.52
2:B:590:SER:C	2:B:591:LYS:HD3	2.34	0.52
2:A:691:GLN:HA	2:A:694:LEU:HD12	1.92	0.51
2:B:726:LEU:HD23	2:B:745:LEU:HD11	1.92	0.51
2:B:873:ASN:O	2:B:878:LYS:HE3	2.10	0.51
2:A:726:LEU:HD23	2:A:745:LEU:HD11	1.92	0.51
2:A:867:LEU:HD13	2:A:881:ILE:HG12	1.93	0.51
2:B:795:HIS:CD2	3:B:903:SO4:O2	2.64	0.51
2:A:825:GLU:HG3	2:A:862:ILE:HG23	1.91	0.51
2:B:826:THR:O	2:B:829:GLN:N	2.40	0.51
2:A:600:THR:C	2:A:602:ASP:H	2.19	0.51
2:A:786:LYS:HB3	2:A:787:PRO:HD3	1.92	0.51
2:A:704:PHE:CZ	2:A:705:ILE:HG13	2.46	0.50
2:A:653:LEU:O	2:A:657:ILE:HG13	2.10	0.50
2:A:709:ILE:HD11	2:A:725:VAL:HG12	1.91	0.50
2:A:793:SER:HB3	2:A:840:VAL:HG21	1.93	0.50
2:B:666:ALA:O	2:B:670:VAL:HG13	2.12	0.50
2:B:847:ILE:HG23	2:B:851:GLN:HG3	1.92	0.50
2:B:854:TYR:CD1	2:B:854:TYR:C	2.89	0.50
2:A:630:ILE:HG22	2:A:634:LYS:HE3	1.94	0.50
2:A:838:ASN:C	2:A:838:ASN:OD1	2.54	0.50
2:B:595:ALA:O	2:B:596:ILE:C	2.55	0.50
2:B:786:LYS:HB3	2:B:787:PRO:HD3	1.93	0.50
1:C:6:A:H2	2:A:654:GLN:HE22	1.58	0.49
2:A:564:TYR:CE2	2:A:565:ILE:HD13	2.47	0.49
2:B:867:LEU:HD13	2:B:881:ILE:HG12	1.94	0.49
2:A:595:ALA:O	2:A:596:ILE:C	2.56	0.49
2:A:624:VAL:HG22	2:A:628:GLN:OE1	2.12	0.49
2:A:568:ILE:HB	2:A:599:GLU:OE1	2.12	0.49
2:A:641:VAL:O	2:A:645:LEU:HG	2.13	0.49
2:B:660:ILE:O	2:B:660:ILE:HG13	2.13	0.49
2:B:753:THR:HG22	2:B:798:GLY:HA2	1.93	0.49
1:C:7:U:N3	2:A:651:ARG:HD2	2.27	0.49
2:A:575:GLN:O	2:A:579:ARG:HB2	2.13	0.49
2:B:825:GLU:O	2:B:825:GLU:HG2	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:849:HIS:HB2	2:B:856:TYR:CD1	2.47	0.49
2:B:630:ILE:HG22	2:B:634:LYS:HE3	1.94	0.49
2:B:600:THR:C	2:B:602:ASP:H	2.20	0.49
2:A:575:GLN:HE22	2:A:579:ARG:NE	1.95	0.49
2:A:753:THR:HG22	2:A:798:GLY:HA2	1.94	0.49
2:B:825:GLU:HG3	2:B:862:ILE:HG23	1.94	0.49
2:B:859:LEU:C	2:B:861:GLU:H	2.21	0.49
2:A:709:ILE:HD11	2:A:725:VAL:CG1	2.42	0.48
2:B:710:SER:HB3	2:B:744:LYS:NZ	2.28	0.48
2:A:585:LEU:HD21	2:A:596:ILE:CD1	2.44	0.48
2:A:826:THR:O	2:A:829:GLN:N	2.44	0.48
2:B:823:GLY:O	2:B:824:GLY:C	2.56	0.48
2:B:614:ASN:O	2:B:618:GLN:HG3	2.12	0.48
2:B:825:GLU:HG3	2:B:862:ILE:CG2	2.43	0.48
2:B:640:PHE:HD2	2:B:672:SER:CB	2.26	0.48
2:A:631:VAL:O	2:A:635:ILE:HG13	2.13	0.48
2:A:659:CYS:O	2:A:661:LYS:HG3	2.13	0.48
2:A:849:HIS:HB2	2:A:856:TYR:CD1	2.49	0.48
2:B:631:VAL:O	2:B:635:ILE:HG13	2.13	0.48
2:B:764:TYR:O	2:B:765:ILE:C	2.57	0.48
2:B:720:ARG:HA	2:B:757:PHE:CE1	2.49	0.47
2:A:847:ILE:HG22	2:A:848:SER:N	2.30	0.47
2:A:847:ILE:HG23	2:A:851:GLN:HG3	1.96	0.47
1:C:7:U:H3	2:A:618:GLN:NE2	2.11	0.47
2:A:575:GLN:HB2	2:A:612:PHE:CD1	2.50	0.47
2:B:779:HIS:NE2	2:B:810:ILE:HD12	2.29	0.47
1:C:7:U:C2	2:A:651:ARG:HD2	2.49	0.47
2:A:710:SER:HB3	2:A:744:LYS:NZ	2.29	0.47
2:B:656:LEU:HD23	2:B:656:LEU:O	2.14	0.47
2:A:823:GLY:O	2:A:824:GLY:O	2.32	0.47
2:B:796:LYS:H	2:B:796:LYS:HG2	1.51	0.46
2:A:569:HIS:ND1	2:A:570:SER:N	2.63	0.46
2:B:847:ILE:O	2:B:850:LYS:N	2.49	0.46
2:A:568:ILE:HG23	2:A:581:LEU:HD22	1.97	0.46
2:B:719:HIS:O	2:B:720:ARG:C	2.58	0.46
2:B:786:LYS:O	2:B:787:PRO:C	2.58	0.46
2:A:575:GLN:O	2:A:575:GLN:CD	2.58	0.46
2:B:676:TYR:HB2	2:B:680:LEU:HD12	1.98	0.46
2:B:837:GLY:O	2:B:838:ASN:C	2.56	0.46
2:A:567:SER:O	2:A:568:ILE:C	2.58	0.46
2:A:676:TYR:CZ	2:B:719:HIS:HD2	2.34	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:854:TYR:CD1	2:A:854:TYR:C	2.93	0.46
2:A:873:ASN:O	2:A:878:LYS:HE3	2.15	0.46
2:A:565:ILE:HG23	2:A:565:ILE:O	2.16	0.46
2:A:794:ILE:O	2:A:836:TYR:CD2	2.63	0.46
2:B:641:VAL:O	2:B:645:LEU:HG	2.15	0.46
2:A:640:PHE:HD2	2:A:672:SER:CB	2.28	0.46
2:A:642:GLU:HG3	2:B:714:ILE:CB	2.38	0.46
2:A:656:LEU:HD23	2:A:656:LEU:O	2.16	0.46
2:A:676:TYR:HB2	2:A:680:LEU:CD1	2.46	0.46
2:B:624:VAL:HG22	2:B:628:GLN:OE1	2.14	0.46
2:B:796:LYS:HE2	2:B:836:TYR:HE2	1.80	0.46
2:A:640:PHE:CE1	2:A:653:LEU:CD1	2.99	0.46
2:A:884:LEU:HD23	2:A:884:LEU:HA	1.68	0.46
2:B:624:VAL:HG22	2:B:628:GLN:HB2	1.98	0.46
2:A:582:GLN:NE2	2:A:615:TYR:HB3	2.26	0.45
2:A:720:ARG:HA	2:A:757:PHE:CE1	2.51	0.45
2:B:709:ILE:HD11	2:B:725:VAL:HG12	1.97	0.45
1:D:7:U:N3	2:B:651:ARG:HD3	2.31	0.45
2:B:847:ILE:HG22	2:B:848:SER:N	2.31	0.45
2:A:642:GLU:OE1	2:B:712:SER:HA	2.16	0.45
2:A:660:ILE:O	2:A:660:ILE:HG13	2.16	0.45
2:B:659:CYS:O	2:B:661:LYS:HG3	2.16	0.45
1:D:7:U:C2	2:B:651:ARG:HD3	2.52	0.45
2:A:624:VAL:HG22	2:A:628:GLN:HB2	1.98	0.45
2:A:642:GLU:OE2	2:B:715:ASP:OD1	2.34	0.45
2:B:615:TYR:O	2:B:619:LYS:HG2	2.17	0.45
2:A:709:ILE:O	2:A:710:SER:C	2.59	0.45
2:A:858:ARG:HG3	2:A:858:ARG:HH11	1.82	0.45
2:A:662:THR:C	2:A:664:GLU:H	2.25	0.45
2:A:754:LEU:HD22	2:A:795:HIS:CD2	2.52	0.45
2:B:703:GLN:OE1	2:B:737:GLN:NE2	2.49	0.45
2:A:764:TYR:O	2:A:765:ILE:C	2.60	0.45
2:A:590:SER:O	2:A:591:LYS:HD3	2.17	0.44
2:A:615:TYR:O	2:A:619:LYS:HG2	2.17	0.44
2:A:755:ASP:HA	2:A:756:PRO:HD3	1.84	0.44
2:A:849:HIS:CE1	2:A:886:LEU:HB2	2.50	0.44
2:A:852:ASN:OD1	2:A:852:ASN:C	2.60	0.44
2:B:706:PHE:CZ	2:B:729:CYS:HB3	2.52	0.44
2:B:709:ILE:HD11	2:B:725:VAL:CG1	2.47	0.44
2:B:769:GLU:OE1	2:B:775:TYR:HA	2.18	0.44
2:A:639:HIS:HD2	2:B:712:SER:HA	1.81	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:596:ILE:H	2:A:596:ILE:HG13	1.59	0.44
2:A:825:GLU:HG3	2:A:862:ILE:CG2	2.46	0.44
2:A:562:ASP:O	2:A:563:GLN:O	2.36	0.44
2:B:709:ILE:O	2:B:710:SER:C	2.58	0.44
2:B:590:SER:O	2:B:591:LYS:HD3	2.18	0.44
2:A:569:HIS:C	2:A:569:HIS:ND1	2.76	0.43
2:A:801:VAL:O	2:A:804:LYS:HB3	2.18	0.43
2:B:600:THR:C	2:B:602:ASP:N	2.76	0.43
2:A:823:GLY:O	2:A:824:GLY:C	2.61	0.43
2:A:716:ILE:HG22	2:A:722:GLY:O	2.19	0.43
2:A:816:ILE:HD13	2:A:848:SER:HB3	1.99	0.43
2:B:884:LEU:HD23	2:B:884:LEU:HA	1.68	0.43
2:A:574:ASP:HB3	2:A:576:HIS:H	1.83	0.43
2:A:837:GLY:O	2:A:838:ASN:C	2.61	0.43
2:A:847:ILE:O	2:A:850:LYS:N	2.52	0.43
2:B:691:GLN:HA	2:B:694:LEU:HD12	2.01	0.43
2:A:642:GLU:CD	2:B:714:ILE:HG22	2.44	0.43
2:A:786:LYS:O	2:A:787:PRO:C	2.61	0.43
2:B:660:ILE:HD11	2:B:697:LEU:HD21	2.01	0.43
2:B:849:HIS:HB2	2:B:856:TYR:CE1	2.54	0.43
2:B:671:ASP:OD1	2:B:671:ASP:N	2.52	0.43
2:A:858:ARG:NH1	2:A:858:ARG:HG3	2.33	0.43
2:A:798:GLY:C	2:A:800:ASN:H	2.25	0.42
2:A:600:THR:C	2:A:602:ASP:N	2.76	0.42
2:A:562:ASP:OD1	2:A:562:ASP:N	2.51	0.42
2:A:660:ILE:HD11	2:A:697:LEU:HD21	2.01	0.42
2:A:565:ILE:HG13	2:A:566:GLY:O	2.18	0.42
2:A:638:PRO:O	2:B:712:SER:HB3	2.19	0.42
2:A:849:HIS:HB2	2:A:856:TYR:CE1	2.55	0.42
2:A:641:VAL:HG21	2:B:715:ASP:HB3	2.02	0.42
2:A:862:ILE:O	2:A:862:ILE:HG22	2.19	0.42
2:B:620:LEU:O	2:B:624:VAL:HB	2.20	0.42
2:B:852:ASN:OD1	2:B:852:ASN:C	2.62	0.42
2:A:639:HIS:CD2	2:B:712:SER:HA	2.54	0.42
2:A:684:LEU:HD12	2:A:684:LEU:O	2.20	0.42
2:B:719:HIS:O	2:B:721:HIS:N	2.53	0.42
2:B:786:LYS:HA	2:B:815:MET:HE1	2.02	0.42
2:A:640:PHE:CD2	2:A:672:SER:HB3	2.52	0.42
1:D:7:U:H3	2:B:618:GLN:HE22	1.66	0.42
2:B:653:LEU:O	2:B:657:ILE:HG13	2.19	0.42
2:B:656:LEU:C	2:B:656:LEU:CD2	2.93	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:801:VAL:O	2:B:804:LYS:HB3	2.20	0.42
2:A:825:GLU:O	2:A:825:GLU:HG2	2.19	0.41
2:A:848:SER:O	2:A:849:HIS:C	2.63	0.41
2:A:580:PHE:CZ	2:A:584:GLN:HG2	2.55	0.41
2:A:639:HIS:C	2:A:643:ILE:HD12	2.45	0.41
2:B:660:ILE:HD11	2:B:697:LEU:CD2	2.50	0.41
2:B:640:PHE:CE1	2:B:653:LEU:CD1	3.03	0.41
1:D:6:A:C6	2:B:651:ARG:NH1	2.88	0.41
2:A:786:LYS:HA	2:A:815:MET:HE1	2.02	0.41
2:A:650:CYS:O	2:A:654:GLN:HG3	2.21	0.41
2:B:798:GLY:C	2:B:800:ASN:H	2.29	0.41
2:A:641:VAL:CG2	2:B:715:ASP:HB3	2.51	0.41
2:A:785:LEU:C	2:A:787:PRO:HD2	2.46	0.41
2:B:684:LEU:HD12	2:B:684:LEU:O	2.21	0.41
2:A:692:LYS:HE2	2:A:692:LYS:HB2	1.77	0.41
2:A:755:ASP:O	2:A:759:ASN:HB3	2.21	0.41
2:B:714:ILE:HD12	2:B:714:ILE:HA	1.94	0.41
2:A:575:GLN:H	2:A:576:HIS:CD2	2.38	0.41
2:B:808:THR:CG2	2:B:811:VAL:HG22	2.33	0.41
2:B:838:ASN:C	2:B:838:ASN:OD1	2.63	0.41
2:B:848:SER:O	2:B:849:HIS:C	2.63	0.41
2:A:569:HIS:HD2	2:A:603:TYR:CE2	2.39	0.41
2:A:582:GLN:HE22	2:A:615:TYR:CB	2.29	0.40
2:B:822:ASN:C	2:B:824:GLY:N	2.78	0.40
2:B:829:GLN:NE2	2:B:833:ASN:ND2	2.68	0.40
2:A:575:GLN:NE2	2:A:579:ARG:HE	1.96	0.40
2:A:620:LEU:O	2:A:624:VAL:HB	2.21	0.40
2:A:798:GLY:C	2:A:800:ASN:N	2.79	0.40
2:A:847:ILE:O	2:A:848:SER:C	2.64	0.40
2:B:785:LEU:O	2:B:786:LYS:C	2.63	0.40
2:B:731:ASP:HA	2:B:768:LYS:HE3	2.03	0.40
2:A:859:LEU:C	2:A:861:GLU:H	2.30	0.40
2:B:804:LYS:HA	2:B:804:LYS:HD2	1.93	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	323/335 (96%)	258 (80%)	51 (16%)	14 (4%)	2	12
2	B	298/335 (89%)	247 (83%)	40 (13%)	11 (4%)	2	15
All	All	621/670 (93%)	505 (81%)	91 (15%)	25 (4%)	2	14

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	563	GLN
2	A	588	LEU
2	A	824	GLY
2	A	853	ASP
2	B	824	GLY
2	B	853	ASP
2	A	568	ILE
2	A	596	ILE
2	B	587	ILE
2	B	596	ILE
2	A	565	ILE
2	A	569	HIS
2	B	588	LEU
2	A	573	LYS
2	A	720	ARG
2	A	860	SER
2	B	799	SER
2	B	860	SER
2	A	595	ALA
2	A	765	ILE
2	B	595	ALA
2	A	601	LYS
2	B	601	LYS
2	B	765	ILE
2	B	589	GLY

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	
2	A	295/303 (97%)	253 (86%)	42 (14%)	3	16
2	B	269/303 (89%)	230 (86%)	39 (14%)	3	15
All	All	564/606 (93%)	483 (86%)	81 (14%)	3	15

All (81) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	562	ASP
2	A	575	GLN
2	A	576	HIS
2	A	621	LEU
2	A	623	GLU
2	A	624	VAL
2	A	626	THR
2	A	637	SER
2	A	657	ILE
2	A	662	THR
2	A	663	ASP
2	A	670	VAL
2	A	684	LEU
2	A	689	VAL
2	A	700	GLU
2	A	702	PHE
2	A	709	ILE
2	A	712	SER
2	A	714	ILE
2	A	715	ASP
2	A	726	LEU
2	A	734	THR
2	A	740	ASN
2	A	748	LEU
2	A	796	LYS
2	A	808	THR
2	A	810	ILE

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Mol	Chain	Res	Type
2	A	828	ILE
2	A	829	GLN
2	A	830	SER
2	A	838	ASN
2	A	842	GLN
2	A	847	ILE
2	A	851	GLN
2	A	854	TYR
2	A	858	ARG
2	A	859	LEU
2	A	871	ILE
2	A	872	ARG
2	A	879	ARG
2	A	884	LEU
2	A	886	LEU
2	B	621	LEU
2	B	623	GLU
2	B	624	VAL
2	B	626	THR
2	B	637	SER
2	B	657	ILE
2	B	663	ASP
2	B	670	VAL
2	B	684	LEU
2	B	689	VAL
2	B	700	GLU
2	B	702	PHE
2	B	709	ILE
2	B	712	SER
2	B	714	ILE
2	B	715	ASP
2	B	726	LEU
2	B	734	THR
2	B	740	ASN
2	B	746	LEU
2	B	748	LEU
2	B	796	LYS
2	B	808	THR
2	B	810	ILE
2	B	828	ILE
2	B	829	GLN
2	B	830	SER

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Mol	Chain	Res	Type
2	B	835	SER
2	B	838	ASN
2	B	842	GLN
2	B	847	ILE
2	B	851	GLN
2	B	854	TYR
2	B	858	ARG
2	B	859	LEU
2	B	871	ILE
2	B	872	ARG
2	B	879	ARG
2	B	884	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	575	GLN
2	A	576	HIS
2	A	582	GLN
2	A	584	GLN
2	A	614	ASN
2	A	618	GLN
2	A	639	HIS
2	A	654	GLN
2	A	679	GLN
2	A	688	HIS
2	A	821	ASN
2	A	829	GLN
2	A	838	ASN
2	B	654	GLN
2	B	679	GLN
2	B	688	HIS
2	B	719	HIS
2	B	773	ASN
2	B	821	ASN
2	B	829	GLN
2	B	838	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	C	7/8 (87%)	0	0
1	D	7/8 (87%)	0	0
All	All	14/16 (87%)	0	0

There are no RNA backbone outliers to report.

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](#)

9 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	$\# Z > 2$
3	SO4	A	905	-	4,4,4	0.23	0	6,6,6	0.20	0
3	SO4	A	906	-	4,4,4	0.26	0	6,6,6	0.13	0
3	SO4	B	909	-	4,4,4	0.26	0	6,6,6	0.22	0
3	SO4	B	908	-	4,4,4	0.24	0	6,6,6	0.17	0
3	SO4	B	903	-	4,4,4	0.27	0	6,6,6	0.35	0
3	SO4	A	902	-	4,4,4	0.40	0	6,6,6	0.26	0
3	SO4	A	904	-	4,4,4	0.27	0	6,6,6	0.13	0
3	SO4	A	901	-	4,4,4	0.25	0	6,6,6	0.26	0
3	SO4	B	907	-	4,4,4	0.20	0	6,6,6	0.34	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.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	905	SO4	1	0
3	B	903	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	8/8 (100%)	2.15	5 (62%) 0 0	40, 49, 89, 90	0
1	D	8/8 (100%)	1.73	4 (50%) 0 0	41, 51, 90, 91	0
2	A	325/335 (97%)	0.80	39 (12%) 9 5	15, 34, 79, 107	0
2	B	300/335 (89%)	0.65	27 (9%) 15 8	14, 34, 70, 106	0
All	All	641/686 (93%)	0.76	75 (11%) 9 5	14, 34, 79, 107	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	586	ASP	7.0
2	A	565	ILE	4.3
2	A	590	SER	4.2
2	B	602	ASP	4.2
2	B	835	SER	3.9
2	A	591	LYS	3.9
2	A	587	ILE	3.8
2	B	833	ASN	3.8
1	C	6	A	3.6
2	B	885	HIS	3.4
2	B	591	LYS	3.4
2	B	603	TYR	3.4
2	B	588	LEU	3.3
2	B	587	ILE	3.3
2	B	875	PRO	3.2
2	A	564	TYR	3.1
2	B	610	ASP	3.1
1	C	5	U	3.1
2	B	609	THR	3.1
2	B	599	GLU	3.1
1	C	8	A	3.0

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Mol	Chain	Res	Type	RSRZ
2	A	584	GLN	3.0
2	B	593	ALA	3.0
2	A	602	ASP	2.9
2	A	659	CYS	2.9
2	A	871	ILE	2.9
2	A	592	ALA	2.9
2	B	868	VAL	2.8
2	A	588	LEU	2.7
1	D	6	A	2.7
2	B	788	ARG	2.7
2	B	870	PRO	2.7
1	D	5	U	2.7
2	A	715	ASP	2.6
2	A	750	ASP	2.6
1	C	7	U	2.6
2	A	649	GLY	2.6
2	A	603	TYR	2.6
2	A	648	HIS	2.5
2	B	613	GLY	2.5
2	A	612	PHE	2.5
2	A	599	GLU	2.5
2	B	853	ASP	2.5
2	A	623	GLU	2.5
2	B	873	ASN	2.5
2	B	601	LYS	2.5
2	A	593	ALA	2.4
2	B	611	SER	2.4
2	A	589	GLY	2.4
2	A	868	VAL	2.4
2	B	648	HIS	2.4
2	A	714	ILE	2.3
2	A	610	ASP	2.3
2	A	566	GLY	2.3
2	B	600	THR	2.3
2	A	595	ALA	2.3
2	B	856	TYR	2.3
2	A	586	ASP	2.3
2	A	608	MET	2.3
2	A	886	LEU	2.3
2	A	879	ARG	2.3
2	A	615	TYR	2.2
2	A	712	SER	2.2

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Mol	Chain	Res	Type	RSRZ
2	B	594	ASP	2.2
2	A	611	SER	2.2
2	A	601	LYS	2.2
2	A	642	GLU	2.1
2	A	722	GLY	2.1
1	C	4	A	2.1
2	B	842	GLN	2.1
1	D	7	U	2.1
2	A	576	HIS	2.1
1	D	1	U	2.1
2	A	582	GLN	2.1
2	A	607	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 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($\text{\AA}^2$)	Q<0.9
3	SO4	B	909	5/5	0.70	0.11	42,44,61,98	0
3	SO4	A	901	5/5	0.86	0.14	13,30,97,111	0
3	SO4	A	904	5/5	0.89	0.11	47,63,80,86	0
3	SO4	B	907	5/5	0.93	0.18	44,44,77,89	0
3	SO4	A	902	5/5	0.93	0.22	8,13,86,96	0
3	SO4	B	903	5/5	0.94	0.18	21,22,50,65	0
3	SO4	A	905	5/5	0.94	0.09	41,41,50,73	0
3	SO4	B	908	5/5	0.94	0.11	43,45,73,79	0
3	SO4	A	906	5/5	0.94	0.09	35,43,69,71	0

6.5 Other polymers [i](#)

There are no such residues in this entry.