



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 06:05 PM UTC

PDB ID : 2BGG / pdb_00002bgg
Title : The structure of a Piwi protein from *Archaeoglobus fulgidus* complexed with a 16nt siRNA duplex.
Authors : Parker, J.S.; Roe, S.M.; Barford, D.
Deposited on : 2004-12-22
Resolution : 2.20 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
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

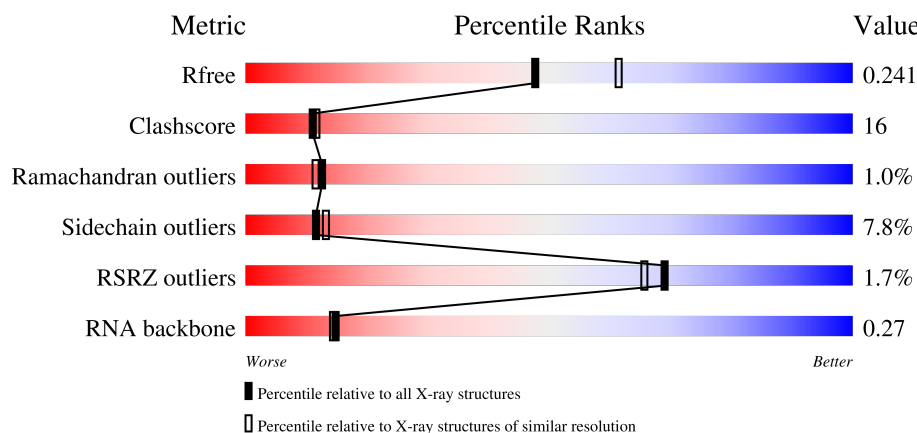
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)
RNA backbone	3983	1052 (2.50-1.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	A	427	
1	B	427	
2	P	8	
2	R	8	

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Mol	Chain	Length	Quality of chain
3	Q	8	12% 38% 25% 25%
3	S	8	12% 25% 12% 12% 38%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7557 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 PROTEIN AF1318.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	395	Total	C	N	O	S	0	2	0
			3237	2097	535	596	9			
1	B	394	Total	C	N	O	S	0	0	0
			3193	2065	530	589	9			

- Molecule 2 is a RNA chain called 5'-R(*UP*UP*CP*GP*AP*CP*GP*CP)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	8	Total	C	N	O	P	0	0	0
			169	75	28	58	8			
2	R	8	Total	C	N	O	P	0	0	0
			169	75	28	58	8			

- Molecule 3 is a RNA chain called 5'-R(*GP*UP*CP*GP*AP*AP*UP*UP)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	Q	8	Total	C	N	O	P	0	0	0
			170	76	29	57	8			
3	S	5	Total	C	N	O	P	0	0	0
			108	48	20	35	5			

- Molecule 4 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mn	0	0
			1	1		
4	B	1	Total	Mn	0	0
			1	1		

- Molecule 5 is water.

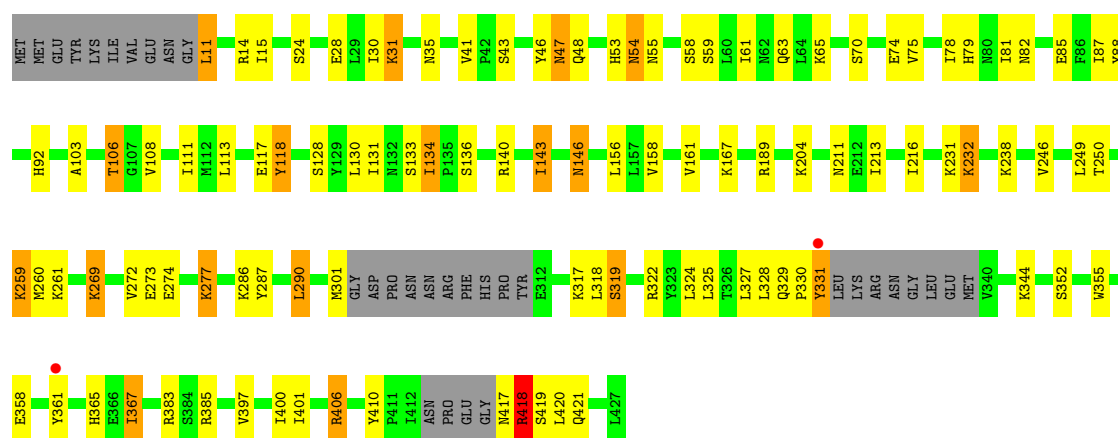
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	283	Total 283	O 283	0	0
5	B	188	Total 188	O 188	0	0
5	P	12	Total 12	O 12	0	0
5	Q	20	Total 20	O 20	0	0
5	R	3	Total 3	O 3	0	0
5	S	3	Total 3	O 3	0	0

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.

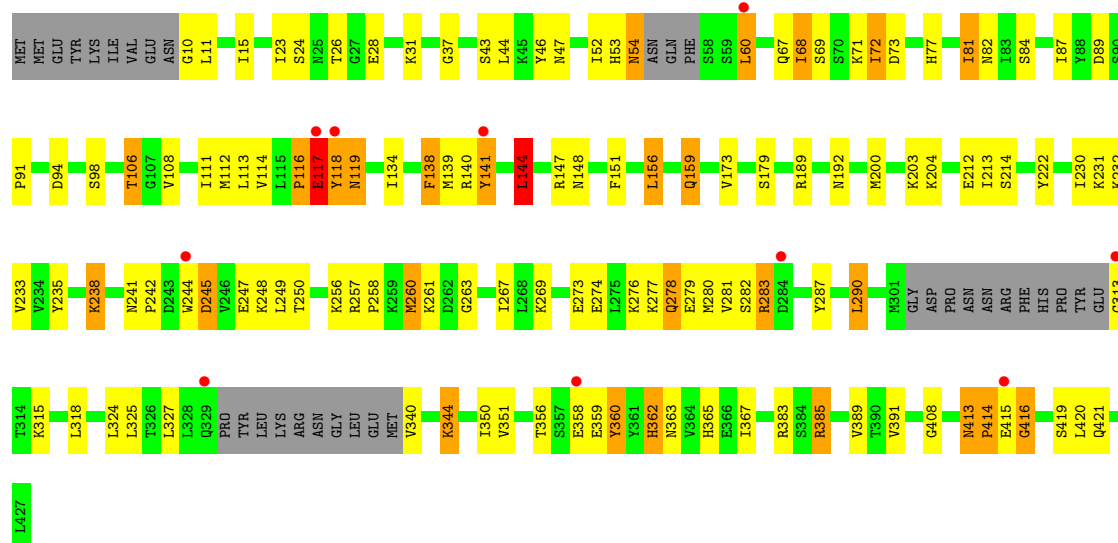
• Molecule 1: PROTEIN AF1318

Chain A: 

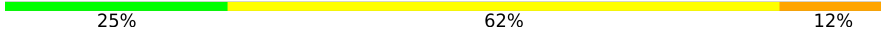


• Molecule 1: PROTEIN AF1318

Chain B: 



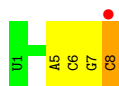
• Molecule 2: 5'-R(*UP*UP*CP*GP*AP*CP*GP*CP)-3'

Chain P:  25% 62% 12%



• Molecule 2: 5'-R(*UP*UP*CP*GP*AP*CP*GP*CP)-3'

Chain R:  12% 50% 38% 12%



• Molecule 3: 5'-R(*GP*UP*CP*GP*AP*AP*UP*UP)-3'

Chain Q:  12% 38% 25% 25%



• Molecule 3: 5'-R(*GP*UP*CP*GP*AP*AP*UP*UP)-3'

Chain S:  12% 12% 25% 12% 12% 38%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	51.92Å 61.21Å 104.06Å 76.54° 76.14° 79.35°	Depositor
Resolution (Å)	99.01 – 2.20 99.01 – 2.20	Depositor EDS
% Data completeness (in resolution range)	96.7 (99.01-2.20) 96.7 (99.01-2.20)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.180 , 0.243 0.180 , 0.241	Depositor DCC
R_{free} test set	2972 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	27.3	Xtriage
Anisotropy	0.360	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 55.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7557	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.03% 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

5.1 Standard geometry

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

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	1.37	10/3316 (0.3%)	1.24	14/4498 (0.3%)
1	B	1.13	5/3269 (0.2%)	1.19	16/4433 (0.4%)
2	P	1.20	0/187	1.70	4/287 (1.4%)
2	R	0.86	0/187	1.29	0/287
3	Q	1.02	1/189 (0.5%)	1.63	5/292 (1.7%)
3	S	0.75	0/120	2.32	6/185 (3.2%)
All	All	1.23	16/7268 (0.2%)	1.27	45/9982 (0.5%)

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
1	A	0	1
1	B	0	3
All	All	0	4

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	161	VAL	CA-CB	8.55	1.64	1.54
1	A	133	SER	CA-C	7.86	1.60	1.52
1	A	30	ILE	CA-CB	7.42	1.63	1.54
1	B	389	VAL	CA-CB	7.18	1.65	1.54
1	A	143	ILE	CA-CB	7.06	1.63	1.54
1	A	397	VAL	CA-CB	6.88	1.62	1.54
1	B	159	GLN	CB-CG	-6.57	1.32	1.52
1	A	216	ILE	CA-CB	6.10	1.61	1.54
1	A	272	VAL	CA-CB	5.70	1.61	1.54
3	Q	11	C	C1'-N1	5.31	1.56	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	167	LYS	C-O	-5.31	1.19	1.24
1	B	52	ILE	CA-CB	5.24	1.60	1.54
1	A	134	ILE	CA-CB	5.23	1.60	1.54
1	A	246	VAL	CA-CB	5.13	1.59	1.53
1	B	413	ASN	CA-C	5.07	1.58	1.52
1	B	351	VAL	CA-CB	5.06	1.60	1.54

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	S	9	G	O3'-P-O5'	-15.11	81.34	104.00
3	S	10	U	O3'-P-O5'	-10.72	87.91	104.00
3	S	10	U	OP1-P-O3'	-9.04	80.89	108.00
3	S	9	G	OP1-P-O3'	-8.37	82.88	108.00
1	B	117	GLU	N-CA-C	8.24	128.36	110.80
2	P	6	C	C5'-C4'-C3'	-7.52	104.72	116.00
3	Q	11	C	C5'-C4'-C3'	-7.43	104.86	116.00
2	P	2	U	N1-C1'-C2'	7.01	122.52	112.00
1	A	59	SER	N-CA-C	6.59	118.12	111.07
1	B	344	LYS	CA-C-N	-6.57	112.24	120.51
1	B	344	LYS	C-N-CA	-6.57	112.24	120.51
1	A	75	VAL	N-CA-C	-6.53	104.65	111.58
1	B	192	ASN	N-CA-C	6.29	119.11	111.82
1	A	325	LEU	CB-CG-CD2	-6.28	91.88	110.70
2	P	1	U	N1-C1'-C2'	-6.13	104.80	114.00
1	B	144	LEU	CA-CB-CG	6.06	137.51	116.30
1	A	41	VAL	CA-C-N	-6.06	113.73	119.85
1	A	41	VAL	C-N-CA	-6.06	113.73	119.85
3	Q	10	U	C3'-C2'-C1'	-5.99	95.31	101.30
1	B	200	MET	CG-SD-CE	-5.94	87.83	100.90
3	Q	13	A	O3'-P-O5'	-5.81	95.28	104.00
2	P	1	U	O5'-C5'-C4'	-5.80	103.00	111.70
1	B	391	VAL	N-CA-C	-5.78	106.84	111.81
3	S	13	A	N9-C1'-C2'	5.72	120.59	112.00
1	B	156	LEU	CA-CB-CG	5.71	136.29	116.30
1	A	397	VAL	CA-C-N	5.70	127.92	120.28
1	A	397	VAL	C-N-CA	5.70	127.92	120.28
1	A	358	GLU	CA-C-N	5.67	128.44	120.28
1	A	358	GLU	C-N-CA	5.67	128.44	120.28
3	S	10	U	OP2-P-O3'	-5.52	91.44	108.00
3	Q	9	G	O4'-C4'-C3'	-5.49	98.51	104.00
1	B	89	ASP	N-CA-C	5.45	119.98	113.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	141	TYR	N-CA-CB	5.44	118.11	110.12
1	B	214	SER	CA-C-N	-5.43	114.11	119.92
1	B	214	SER	C-N-CA	-5.43	114.11	119.92
1	B	119	ASN	N-CA-C	5.36	122.23	110.80
1	A	158	VAL	N-CA-CB	5.34	116.43	110.51
1	B	159	GLN	CB-CA-C	-5.32	102.49	110.90
1	A	35	ASN	N-CA-C	5.32	116.88	111.14
1	B	68	ILE	N-CA-C	-5.31	105.32	110.42
3	Q	9	G	C4'-C3'-C2'	-5.30	97.30	102.60
1	A	117	GLU	N-CA-C	5.28	122.05	110.80
1	A	365	HIS	N-CA-C	5.09	117.22	111.11
1	B	94	ASP	N-CA-C	5.07	116.80	111.28
1	A	74	GLU	N-CA-C	5.04	119.50	112.90

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	259	LYS	Peptide
1	B	116	PRO	Peptide
1	B	117	GLU	Peptide
1	B	414	PRO	Peptide

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	3237	0	3255	90	0
1	B	3193	0	3222	123	0
2	P	169	0	87	8	0
2	R	169	0	87	1	0
3	Q	170	0	86	12	0
3	S	108	0	55	7	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	283	0	0	23	0
5	B	188	0	0	37	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	P	12	0	0	2	0
5	Q	20	0	0	3	0
5	R	3	0	0	0	0
5	S	3	0	0	0	0
All	All	7557	0	6792	226	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 16.

All (226) 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:331:TYR:CE1	2:P:5:A:H5''	1.77	1.18
1:B:385:ARG:HG3	1:B:385:ARG:HH11	1.03	1.17
1:B:282:SER:HB2	5:B:2121:HOH:O	1.44	1.13
1:B:421:GLN:HG2	5:B:2094:HOH:O	1.48	1.12
1:A:63:GLN:HE22	1:A:146:ASN:ND2	1.47	1.12
1:B:419:SER:HA	5:B:2181:HOH:O	1.57	1.03
1:B:54:ASN:HB2	5:B:2030:HOH:O	1.58	1.02
3:Q:15:U:H2'	5:Q:2020:HOH:O	1.57	1.02
1:A:78:ILE:HG22	1:A:81:ILE:HD11	1.43	1.01
1:A:329:GLN:NE2	1:A:344:LYS:HB3	1.76	1.00
1:A:63:GLN:NE2	1:A:146:ASN:HD21	1.59	1.00
1:B:385:ARG:HH11	1:B:385:ARG:CG	1.78	0.97
1:A:78:ILE:CG2	1:A:81:ILE:HD11	1.95	0.95
3:Q:11:C:H5''	3:Q:11:C:H6	1.31	0.93
1:A:14:ARG:HG2	5:A:2011:HOH:O	1.71	0.91
1:A:82[B]:ASN:OD1	1:B:416:GLY:HA2	1.71	0.91
1:B:118:TYR:H	1:B:140:ARG:HH11	1.16	0.88
1:B:11:LEU:HD12	1:B:23:ILE:HG21	1.55	0.86
3:Q:11:C:H5''	3:Q:11:C:C6	2.09	0.86
1:A:261:LYS:HA	5:A:2191:HOH:O	1.75	0.86
1:A:63:GLN:HE22	1:A:146:ASN:HD21	0.89	0.85
1:B:414:PRO:HD2	5:B:2002:HOH:O	1.75	0.85
1:B:313:GLY:HA2	5:B:2135:HOH:O	1.75	0.84
1:B:385:ARG:HG3	1:B:385:ARG:NH1	1.77	0.84
1:B:231:LYS:NZ	1:B:274:GLU:OE1	2.15	0.80
1:A:383:ARG:NH2	3:Q:16:U:O2'	2.14	0.80
1:B:416:GLY:O	5:B:2176:HOH:O	2.00	0.79
1:B:235:TYR:HD1	1:B:280:MET:HE2	1.48	0.77
1:B:116:PRO:HD2	1:B:117:GLU:OE2	1.83	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:28:GLU:OE1	1:A:31:LYS:NZ	2.19	0.76
1:B:118:TYR:N	1:B:140:ARG:HH11	1.84	0.75
1:A:106:THR:HG21	5:A:2043:HOH:O	1.86	0.74
1:A:331:TYR:CE1	2:P:5:A:C5'	2.66	0.74
1:A:329:GLN:HE22	1:A:344:LYS:HB3	1.53	0.73
1:A:418:ARG:HG3	1:A:418:ARG:HH11	1.54	0.73
1:B:241:ASN:HA	5:B:2107:HOH:O	1.88	0.73
1:A:232:LYS:NZ	1:A:232:LYS:HB3	2.04	0.72
1:B:242:PRO:CD	5:B:2107:HOH:O	2.37	0.72
1:B:173:VAL:HG21	1:B:204:LYS:HG3	1.71	0.72
1:B:106:THR:HG21	5:B:2048:HOH:O	1.89	0.71
3:Q:16:U:C6	5:Q:2017:HOH:O	2.44	0.71
1:B:118:TYR:HA	1:B:140:ARG:NH1	2.05	0.70
1:A:406:ARG:HH11	1:A:406:ARG:HG3	1.56	0.70
1:B:118:TYR:H	1:B:140:ARG:NH1	1.89	0.70
1:A:63:GLN:NE2	1:A:146:ASN:ND2	2.28	0.69
1:B:325:LEU:HD23	5:B:2008:HOH:O	1.94	0.68
1:A:82[B]:ASN:OD1	1:B:416:GLY:CA	2.41	0.68
1:A:78:ILE:HG22	1:A:81:ILE:CD1	2.20	0.68
1:A:128:SER:HB2	1:A:410:TYR:O	1.94	0.68
1:B:383:ARG:NH2	3:S:12:G:H21	1.90	0.67
1:A:231:LYS:HE3	1:A:274:GLU:OE1	1.94	0.67
1:A:331:TYR:CZ	2:P:5:A:H5''	2.28	0.67
2:P:4:G:N3	5:P:2008:HOH:O	2.27	0.67
1:A:261:LYS:HE3	5:A:2236:HOH:O	1.95	0.66
1:B:325:LEU:HG	5:B:2130:HOH:O	1.95	0.66
1:B:274:GLU:HA	1:B:277:LYS:HE3	1.76	0.66
1:A:232:LYS:HB3	1:A:232:LYS:HZ2	1.62	0.65
1:B:11:LEU:CD1	1:B:23:ILE:HG21	2.27	0.64
1:A:406:ARG:HG3	1:A:406:ARG:NH1	2.11	0.64
1:B:269:LYS:O	1:B:273:GLU:HG2	1.96	0.63
1:A:329:GLN:OE1	1:A:330:PRO:HD2	1.98	0.63
1:B:235:TYR:HA	1:B:280:MET:HE1	1.80	0.63
1:A:106:THR:CG2	5:A:2043:HOH:O	2.45	0.63
1:A:269:LYS:HE2	1:A:352:SER:OG	1.98	0.63
1:B:54:ASN:C	1:B:54:ASN:HD22	2.08	0.61
1:A:231:LYS:CE	1:A:274:GLU:OE1	2.49	0.61
1:A:111:ILE:O	1:A:136:SER:HA	2.01	0.60
1:B:242:PRO:HD2	5:B:2107:HOH:O	1.99	0.60
3:S:10:U:OP2	3:S:10:U:H6	1.84	0.60
1:B:203:LYS:HE3	1:B:244:TRP:CH2	2.36	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:273:GLU:O	1:A:277:LYS:HD2	2.03	0.59
1:B:118:TYR:CD2	1:B:119:ASN:N	2.71	0.59
1:B:44:LEU:CD1	1:B:81:ILE:HD11	2.33	0.58
1:A:211:ASN:OD1	5:A:2144:HOH:O	2.17	0.58
1:B:414:PRO:O	5:B:2175:HOH:O	2.16	0.57
1:A:11:LEU:N	5:A:2007:HOH:O	2.36	0.57
1:A:355:TRP:CZ2	1:A:367:ILE:HD12	2.39	0.57
1:B:60:LEU:HD11	1:B:144:LEU:HD22	1.87	0.57
1:A:55:ASN:HB2	1:A:61:ILE:CD1	2.34	0.57
1:B:117:GLU:HG3	1:B:119:ASN:HB2	1.86	0.56
1:A:211:ASN:HB2	5:A:2144:HOH:O	2.03	0.56
1:B:118:TYR:CA	1:B:140:ARG:HH11	2.19	0.56
1:A:54:ASN:HB3	1:A:88:TYR:CE1	2.40	0.56
1:B:118:TYR:CA	1:B:140:ARG:NH1	2.69	0.56
1:A:204:LYS:CG	5:A:2140:HOH:O	2.54	0.55
1:B:360:TYR:HE2	5:B:2133:HOH:O	1.88	0.55
1:B:383:ARG:HH11	3:S:13:A:H1'	1.72	0.54
1:A:204:LYS:HG2	5:A:2140:HOH:O	2.07	0.54
1:B:212:GLU:CD	5:B:2094:HOH:O	2.50	0.54
1:A:319:SER:HB3	1:A:322:ARG:NH1	2.23	0.54
1:A:400:ILE:HD11	5:A:2104:HOH:O	2.07	0.54
1:B:212:GLU:HG3	1:B:232:LYS:HB3	1.90	0.54
1:B:383:ARG:HH11	3:S:13:A:C1'	2.21	0.54
1:B:118:TYR:HD2	1:B:119:ASN:H	1.56	0.53
1:B:44:LEU:HD12	1:B:81:ILE:HD11	1.89	0.53
1:B:116:PRO:HA	1:B:141:TYR:HD1	1.74	0.53
1:B:141:TYR:HA	1:B:144:LEU:HD13	1.91	0.53
1:B:118:TYR:N	1:B:140:ARG:NH1	2.52	0.53
1:B:203:LYS:HE3	1:B:244:TRP:HH2	1.73	0.53
1:A:331:TYR:CD1	2:P:5:A:H5''	2.40	0.52
1:B:73:ASP:OD2	5:B:2038:HOH:O	2.18	0.52
1:B:189:ARG:HG3	5:B:2084:HOH:O	2.08	0.52
1:B:258:PRO:HD3	5:B:2084:HOH:O	2.09	0.52
3:Q:10:U:H2'	3:Q:11:C:O4'	2.08	0.52
3:S:10:U:OP2	3:S:10:U:C6	2.63	0.52
1:A:269:LYS:NZ	5:A:2195:HOH:O	2.43	0.52
1:B:340:VAL:O	1:B:340:VAL:HG23	2.10	0.51
5:A:2021:HOH:O	3:Q:14:A:H5''	2.10	0.51
1:B:383:ARG:HD2	5:B:2160:HOH:O	2.09	0.51
1:A:355:TRP:HZ2	1:A:367:ILE:HD12	1.76	0.51
1:B:238:LYS:HB2	1:B:280:MET:HE1	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:54:ASN:CB	5:B:2030:HOH:O	2.35	0.51
1:A:406:ARG:HH11	1:A:406:ARG:CG	2.23	0.50
1:A:24:SER:O	3:Q:16:U:H5'	2.12	0.50
1:B:249:LEU:O	1:B:287:TYR:HA	2.11	0.50
1:B:54:ASN:OD1	5:B:2032:HOH:O	2.19	0.50
1:A:63:GLN:CD	1:A:146:ASN:HD21	2.19	0.50
1:B:263:GLY:O	1:B:267:ILE:HG12	2.12	0.50
1:B:10:GLY:HA2	5:B:2004:HOH:O	2.11	0.49
1:B:235:TYR:HA	1:B:280:MET:CE	2.42	0.49
1:B:138:PHE:CD1	1:B:138:PHE:N	2.81	0.49
1:A:331:TYR:CD1	2:P:5:A:C5'	2.96	0.49
1:B:82:ASN:HB3	5:B:2029:HOH:O	2.12	0.49
1:B:281:VAL:HA	5:B:2106:HOH:O	2.12	0.49
1:B:415:GLU:O	1:B:416:GLY:C	2.55	0.49
1:A:259:LYS:HD2	5:A:2183:HOH:O	2.13	0.48
1:B:327:LEU:HD21	5:B:2008:HOH:O	2.13	0.48
1:A:65:LYS:NZ	1:A:85:GLU:HG3	2.27	0.48
1:B:117:GLU:HG3	1:B:119:ASN:CB	2.44	0.48
1:B:116:PRO:HD2	1:B:117:GLU:CD	2.37	0.48
1:A:301:MET:C	5:A:2223:HOH:O	2.57	0.48
1:A:48:GLN:HG2	1:A:108:VAL:HG12	1.96	0.48
1:B:117:GLU:HB3	1:B:118:TYR:C	2.39	0.47
1:A:317:LYS:NZ	1:A:361[B]:TYR:OH	2.48	0.47
1:B:179:SER:OG	1:B:250:THR:OG1	2.28	0.47
1:B:315:LYS:HE3	1:B:365:HIS:CE1	2.50	0.47
1:A:65:LYS:HZ2	1:A:85:GLU:HG3	1.79	0.47
1:A:331:TYR:C	3:Q:16:U:O2	2.57	0.47
1:B:53:HIS:O	1:B:87:ILE:HA	2.14	0.47
3:Q:10:U:H6	3:Q:10:U:O5'	1.96	0.47
1:B:383:ARG:HH22	3:S:12:G:H21	1.60	0.47
2:P:6:C:H2'	2:P:7:G:C8	2.50	0.47
1:A:14:ARG:CG	5:A:2011:HOH:O	2.42	0.47
1:A:213:ILE:HD12	5:A:2147:HOH:O	2.13	0.47
1:B:242:PRO:HD3	5:B:2107:HOH:O	2.06	0.47
1:B:138:PHE:N	1:B:138:PHE:HD1	2.12	0.46
1:B:81:ILE:HG22	5:B:2037:HOH:O	2.16	0.46
1:A:46:TYR:O	1:A:47:ASN:HB2	2.15	0.46
1:A:250:THR:CG2	1:A:290:LEU:HD22	2.46	0.46
1:B:274:GLU:O	1:B:278:GLN:HG2	2.15	0.46
1:A:92:HIS:HB2	5:A:2087:HOH:O	2.16	0.46
1:A:417:ASN:C	1:A:419:SER:H	2.23	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:383:ARG:NH1	3:S:13:A:H1'	2.31	0.46
1:B:112:MET:HB3	1:B:139:MET:HE2	1.98	0.46
1:B:362:HIS:HD2	1:B:363:ASN:OD1	1.99	0.46
1:A:130:LEU:CD2	1:A:134:ILE:HG22	2.45	0.46
1:B:28:GLU:HG2	5:B:2021:HOH:O	2.15	0.46
1:A:331:TYR:CD1	1:A:331:TYR:N	2.84	0.45
1:A:249:LEU:O	1:A:287:TYR:HA	2.16	0.45
1:A:418:ARG:O	1:A:418:ARG:HD3	2.17	0.45
1:A:189:ARG:HD3	5:A:2137:HOH:O	2.17	0.45
1:B:60:LEU:CD1	1:B:144:LEU:HD22	2.46	0.45
1:B:245:ASP:CB	5:B:2109:HOH:O	2.64	0.45
1:B:245:ASP:HA	5:B:2109:HOH:O	2.15	0.45
1:A:417:ASN:O	1:A:419:SER:N	2.50	0.45
1:B:173:VAL:CG2	1:B:204:LYS:HG3	2.44	0.45
1:B:362:HIS:CD2	1:B:363:ASN:OD1	2.70	0.45
3:Q:11:C:H5'	5:Q:2008:HOH:O	2.17	0.45
1:B:23:ILE:O	1:B:24:SER:HB2	2.18	0.44
1:A:287:TYR:CE1	5:A:2213:HOH:O	2.69	0.44
2:R:7:G:C6	2:R:8:C:N4	2.86	0.44
1:B:203:LYS:HG2	1:B:244:TRP:CH2	2.53	0.44
1:B:278:GLN:O	1:B:279:GLU:HB2	2.17	0.43
1:A:54:ASN:ND2	5:A:2046:HOH:O	2.50	0.43
1:B:37:GLY:HA2	1:B:77:HIS:CD2	2.53	0.43
1:B:385:ARG:HA	1:B:385:ARG:HD2	1.66	0.43
1:A:213:ILE:O	1:A:213:ILE:HG13	2.18	0.43
1:B:235:TYR:CD1	1:B:280:MET:HE2	2.39	0.43
1:B:385:ARG:CG	1:B:385:ARG:NH1	2.49	0.43
1:B:260:MET:HE1	5:B:2099:HOH:O	2.19	0.43
1:A:286:LYS:HA	1:A:286:LYS:HD3	1.90	0.42
1:B:68:ILE:HG22	1:B:72:ILE:HD12	2.01	0.42
1:B:278:GLN:HA	5:B:2119:HOH:O	2.18	0.42
1:A:55:ASN:HB3	1:A:58:SER:HB2	2.01	0.42
1:A:78:ILE:HG22	1:A:81:ILE:CG1	2.48	0.42
1:B:276:LYS:HG2	1:B:283:ARG:HA	2.01	0.42
1:A:79:HIS:HE1	5:A:2066:HOH:O	2.02	0.42
1:A:118:TYR:HB2	1:A:140:ARG:NH2	2.35	0.42
1:A:213:ILE:HD13	1:A:401:ILE:HD11	2.02	0.42
1:B:222:TYR:OH	1:B:257:ARG:HD2	2.20	0.42
1:A:417:ASN:C	1:A:419:SER:N	2.78	0.42
1:B:408:GLY:HA2	5:B:2169:HOH:O	2.18	0.42
1:B:414:PRO:HG2	5:B:2093:HOH:O	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:269:LYS:HB2	1:A:269:LYS:HE3	1.73	0.42
1:A:319:SER:HB3	1:A:322:ARG:HH11	1.84	0.42
1:B:60:LEU:CD1	1:B:144:LEU:CD2	2.98	0.41
1:B:148:ASN:ND2	1:B:151:PHE:HD1	2.18	0.41
3:Q:11:C:C6	3:Q:11:C:C5'	2.94	0.41
1:B:46:TYR:O	1:B:47:ASN:HB2	2.20	0.41
1:B:114:VAL:HG12	1:B:141:TYR:CE1	2.56	0.41
1:A:418:ARG:HG3	1:A:418:ARG:NH1	2.27	0.41
1:B:230:ILE:O	1:B:233:VAL:HG12	2.20	0.41
1:A:417:ASN:HB2	5:A:2274:HOH:O	2.20	0.41
1:B:356:THR:O	1:B:359:GLU:HB3	2.20	0.41
1:A:259:LYS:HD2	1:A:259:LYS:HA	1.93	0.41
1:B:54:ASN:C	1:B:54:ASN:ND2	2.78	0.41
1:A:318:LEU:HD11	1:A:324:LEU:HG	2.02	0.41
1:B:10:GLY:N	5:B:2003:HOH:O	2.53	0.41
1:B:67:GLN:OE1	1:B:71:LYS:HE2	2.20	0.41
1:B:91:PRO:HG3	5:B:2032:HOH:O	2.20	0.41
1:B:117:GLU:OE1	1:B:117:GLU:HA	2.19	0.41
1:B:420:LEU:HD23	1:B:420:LEU:HA	1.91	0.41
1:B:189:ARG:HD2	1:B:256:LYS:HE3	2.03	0.41
1:A:53:HIS:O	1:A:87:ILE:HA	2.21	0.40
1:B:118:TYR:HD2	1:B:119:ASN:N	2.12	0.40
1:A:103:ALA:O	1:A:106:THR:HB	2.21	0.40
1:A:143:ILE:HD13	5:P:2002:HOH:O	2.20	0.40
1:B:318:LEU:HD11	1:B:324:LEU:HG	2.03	0.40
2:P:5:A:H2'	2:P:6:C:C6	2.56	0.40
1:B:111:ILE:HD12	1:B:134:ILE:HG21	2.04	0.40
1:B:280:MET:HE3	1:B:280:MET:HB3	1.94	0.40
1:A:54:ASN:HB3	1:A:88:TYR:CD1	2.57	0.40
1:B:413:ASN:ND2	1:B:416:GLY:H	2.20	0.40
1:A:131:ILE:O	1:A:420:LEU:HD21	2.21	0.40
1:B:108:VAL:HG23	1:B:134:ILE:HD11	2.02	0.40
1:B:290:LEU:HD12	1:B:350:ILE:HA	2.02	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	389/427 (91%)	373 (96%)	12 (3%)	4 (1%)	12	11
1	B	386/427 (90%)	368 (95%)	14 (4%)	4 (1%)	12	11
All	All	775/854 (91%)	741 (96%)	26 (3%)	8 (1%)	12	11

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	118	TYR
1	A	260	MET
1	A	418	ARG
1	B	118	TYR
1	B	360	TYR
1	A	15	ILE
1	B	15	ILE
1	B	416	GLY

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	
1	A	368/394 (93%)	344 (94%)	24 (6%)	15	18
1	B	363/394 (92%)	330 (91%)	33 (9%)	9	9
All	All	731/788 (93%)	674 (92%)	57 (8%)	11	13

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	LEU
1	A	31	LYS
1	A	43	SER
1	A	47	ASN
1	A	54	ASN
1	A	70	SER
1	A	106	THR
1	A	113	LEU
1	A	146	ASN
1	A	156	LEU
1	A	232	LYS
1	A	238	LYS
1	A	269	LYS
1	A	277	LYS
1	A	290	LEU
1	A	319	SER
1	A	327	LEU
1	A	328	LEU
1	A	331	TYR
1	A	367	ILE
1	A	385	ARG
1	A	406	ARG
1	A	418	ARG
1	A	421	GLN
1	B	26	THR
1	B	31	LYS
1	B	43	SER
1	B	54	ASN
1	B	60	LEU
1	B	69	SER
1	B	72	ILE
1	B	81	ILE
1	B	84	SER
1	B	98	SER
1	B	106	THR
1	B	113	LEU
1	B	117	GLU
1	B	138	PHE
1	B	144	LEU
1	B	147	ARG
1	B	156	LEU
1	B	159	GLN
1	B	213	ILE

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Mol	Chain	Res	Type
1	B	238	LYS
1	B	245	ASP
1	B	247	GLU
1	B	248	LYS
1	B	260	MET
1	B	261	LYS
1	B	278	GLN
1	B	283	ARG
1	B	290	LEU
1	B	344	LYS
1	B	358	GLU
1	B	362	HIS
1	B	367	ILE
1	B	385	ARG

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

Mol	Chain	Res	Type
1	A	54	ASN
1	A	99	GLN
1	A	146	ASN
1	A	148	ASN
1	A	172	ASN
1	B	54	ASN
1	B	79	HIS
1	B	278	GLN
1	B	329	GLN
1	B	362	HIS
1	B	403	ASN
1	B	423	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	P	7/8 (87%)	0	0
2	R	7/8 (87%)	3 (42%)	0
3	Q	8/8 (100%)	3 (37%)	1 (12%)
3	S	4/8 (50%)	1 (25%)	0
All	All	26/32 (81%)	7 (26%)	1 (3%)

All (7) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
3	Q	10	U
3	Q	11	C
3	Q	16	U
2	R	5	A
2	R	6	C
2	R	8	C
3	S	10	U

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
3	Q	9	G

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 2 ligands modelled in this entry, 2 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 [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	395/427 (92%)	-0.34	2 (0%) 87 85	11, 22, 40, 63	2 (0%)
1	B	394/427 (92%)	0.12	10 (2%) 58 55	18, 32, 55, 65	0
2	P	8/8 (100%)	-0.14	0 100 100	20, 35, 78, 86	0
2	R	8/8 (100%)	0.44	1 (12%) 8 6	34, 66, 93, 94	0
3	Q	8/8 (100%)	0.35	0 100 100	42, 47, 78, 87	0
3	S	5/8 (62%)	1.55	1 (20%) 3 2	78, 80, 86, 89	0
All	All	818/886 (92%)	-0.09	14 (1%) 69 66	11, 27, 54, 94	2 (0%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	118	TYR	5.6
1	A	331	TYR	5.6
1	B	141	TYR	3.9
1	A	361[A]	TYR	3.4
1	B	244	TRP	3.1
1	B	117	GLU	2.9
1	B	313	GLY	2.8
3	S	13	A	2.7
1	B	358	GLU	2.5
2	R	8	C	2.5
1	B	284	ASP	2.4
1	B	415	GLU	2.1
1	B	329	GLN	2.1
1	B	60	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
4	MN	A	1428	1/1	0.99	0.02	16,16,16,16	0
4	MN	B	1428	1/1	1.00	0.02	23,23,23,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.