



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 06:35 AM UTC

PDB ID : 7FO8 / pdb_00007fo8
Title : PanDDA analysis group deposition – Aar2/RNaseH in complex with fragment P07H06 from the F2X-Universal Library
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Deposited on : 2022-08-26
Resolution : 1.72 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

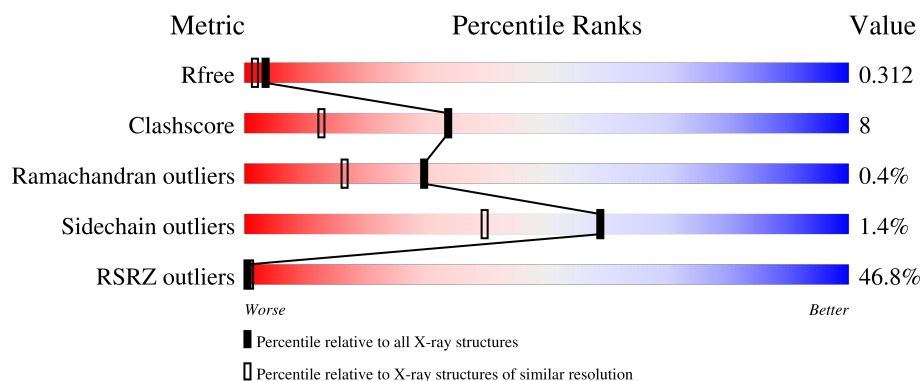
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.72 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	1039 (1.72-1.72)
Clashscore	190562	1049 (1.72-1.72)
Ramachandran outliers	187476	1041 (1.72-1.72)
Sidechain outliers	187428	1041 (1.72-1.72)
RSRZ outliers	180081	1039 (1.72-1.72)

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	258	42% 75% 15% 8%
2	B	308	46% 85% 12%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9205 atoms, of which 4524 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 Pre-mRNA-splicing factor 8.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	237	4068	1287	2060	336	373	12	18	21	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1833	GLY	-	expression tag	UNP P33334
A	1834	ALA	-	expression tag	UNP P33334
A	1835	MET	-	expression tag	UNP P33334

- Molecule 2 is a protein called A1 cistron-splicing factor AAR2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
2	B	300	5044	1654	2464	421	485	20	0	17	0

There are 20 discrepancies between the modelled and reference sequences:

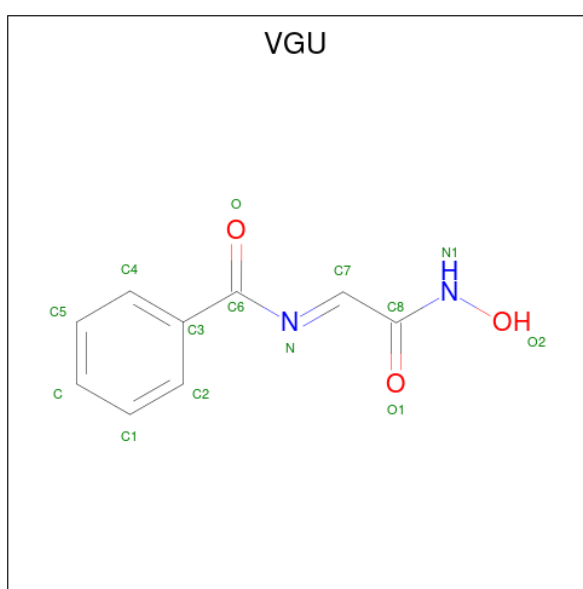
Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	GLY	-	expression tag	UNP P32357
B	-2	ALA	-	expression tag	UNP P32357
B	-1	MET	-	expression tag	UNP P32357
B	0	ALA	-	expression tag	UNP P32357
B	166	SER	LEU	conflict	UNP P32357
B	167	SER	LYS	conflict	UNP P32357
B	?	-	LEU	deletion	UNP P32357
B	?	-	GLN	deletion	UNP P32357
B	?	-	LYS	deletion	UNP P32357
B	?	-	ALA	deletion	UNP P32357
B	?	-	GLY	deletion	UNP P32357
B	?	-	SER	deletion	UNP P32357
B	?	-	LYS	deletion	UNP P32357

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Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	MET	deletion	UNP P32357
B	?	-	GLU	deletion	UNP P32357
B	?	-	ALA	deletion	UNP P32357
B	?	-	LYS	deletion	UNP P32357
B	?	-	ASN	deletion	UNP P32357
B	?	-	GLU	deletion	UNP P32357
B	170	SER	ASP	conflict	UNP P32357

- Molecule 3 is N-[(1E)-2-(hydroxyamino)-2-oxoethylidene]benzamide (CCD ID: VGU) (formula: C₉H₈N₂O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	9	2	3		

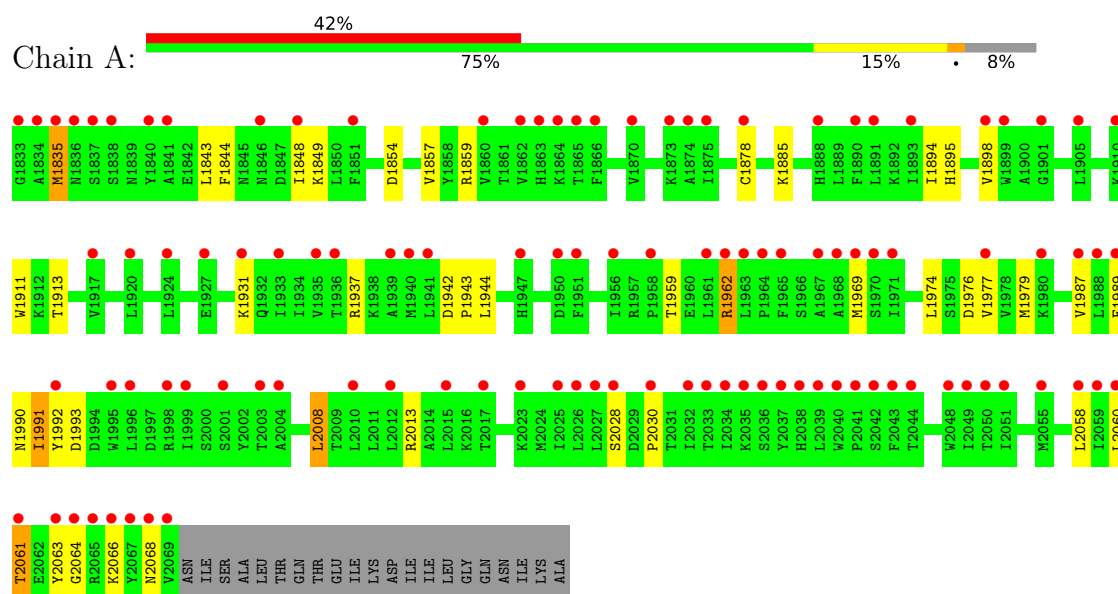
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	35	Total	O	0	0
			35	35		
4	B	44	Total	O	0	0
			44	44		

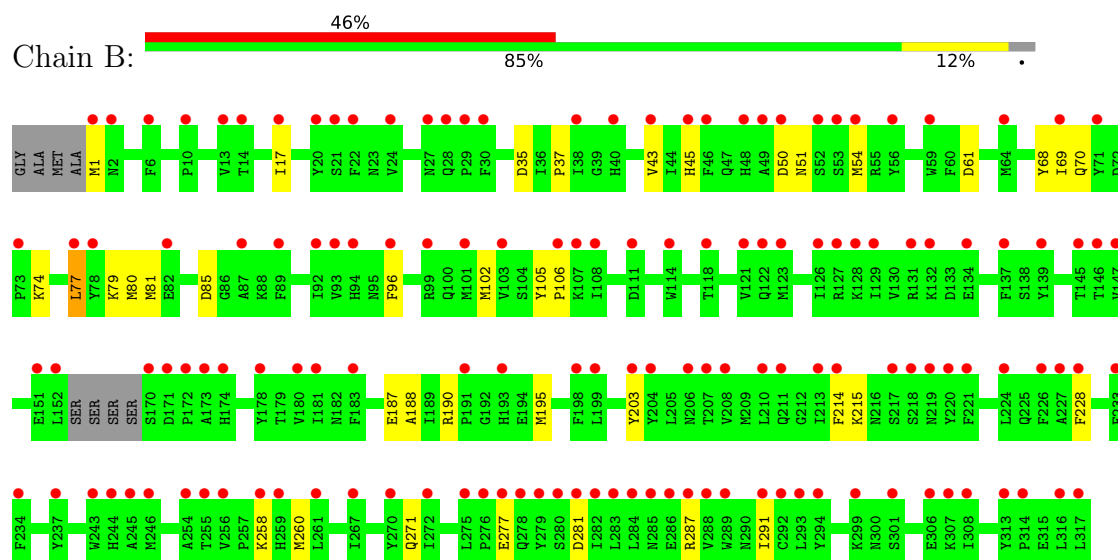
3 Residue-property plots

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: Pre-mRNA-splicing factor 8



• Molecule 2: A1 cistron-splicing factor AAR2



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	88.44Å 81.76Å 92.76Å 90.00° 107.98° 90.00°	Depositor
Resolution (Å)	44.11 – 1.72 44.11 – 1.72	Depositor EDS
% Data completeness (in resolution range)	99.7 (44.11-1.72) 99.9 (44.11-1.72)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.97 (at 1.72Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.265 , 0.314 0.275 , 0.312	Depositor DCC
R_{free} test set	2098 reflections (3.14%)	wwPDB-VP
Wilson B-factor (Å ²)	42.8	Xtriage
Anisotropy	0.381	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 53.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9205	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.04% 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: VGU

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	0.69	0/2149	0.84	3/2911 (0.1%)
2	B	0.69	0/2739	0.83	0/3699
All	All	0.69	0/4888	0.83	3/6610 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	2008	LEU	CB-CG-CD1	-5.71	93.57	110.70
1	A	1990	ASN	N-CA-CB	5.21	118.96	110.37
1	A	1990	ASN	N-CA-C	-5.07	99.54	108.20

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	2008	2060	1974	34	0
2	B	2580	2464	2398	36	0
3	A	14	0	0	0	0
4	A	35	0	0	0	0
4	B	44	0	0	4	0
All	All	4681	4524	4372	69	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 8.

All (69) 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:1974:LEU:CD2	1:A:1989:PHE:CE1	2.46	0.97
1:A:1974:LEU:HD23	1:A:1989:PHE:CE1	1.98	0.96
1:A:1974:LEU:CD2	1:A:1989:PHE:CD1	2.48	0.96
1:A:1962:ARG:O	1:A:2013:ARG:NH1	2.11	0.83
1:A:1974:LEU:HD22	1:A:1989:PHE:CD1	2.24	0.72
1:A:1974:LEU:HD22	1:A:1989:PHE:CE1	2.29	0.68
2:B:77:LEU:HD21	2:B:79:LYS:HE3	1.75	0.67
2:B:50:ASP:OD1	2:B:51:ASN:N	2.26	0.64
1:A:2064:GLY:O	1:A:2068:ASN:N	2.30	0.64
1:A:1974:LEU:HD21	1:A:1989:PHE:CD1	2.34	0.63
2:B:70:GLN:HB3	2:B:81:MET:CE	2.30	0.62
1:A:1895:HIS:O	1:A:1898[A]:VAL:HG22	2.04	0.58
2:B:1:MET:HB3	2:B:35:ASP:HA	1.86	0.57
1:A:1979[B]:MET:HA	1:A:1979[B]:MET:HE2	1.87	0.57
1:A:1977:VAL:HG21	1:A:1987:VAL:HG21	1.86	0.57
2:B:43:VAL:HG13	2:B:43:VAL:O	2.03	0.57
2:B:37:PRO:HD3	2:B:105:TYR:CD1	2.41	0.56
2:B:287:ARG:O	2:B:291:ILE:HD13	2.06	0.56
1:A:1843:LEU:HA	1:A:1849:LYS:HD2	1.89	0.55
1:A:1848:ILE:H	1:A:1931[A]:LYS:HZ2	1.53	0.55
1:A:2061:THR:O	1:A:2064:GLY:N	2.40	0.55
1:A:1878:CYS:SG	1:A:1969:MET:HE3	2.48	0.54
2:B:1:MET:N	4:B:401:HOH:O	2.39	0.53
2:B:277:GLU:CD	2:B:277:GLU:H	2.17	0.51
1:A:1911:TRP:CD1	2:B:195:MET:HE3	2.46	0.51
2:B:70:GLN:HB3	2:B:81:MET:HE1	1.93	0.50
2:B:96:PHE:HB2	2:B:102:MET:HE3	1.93	0.50
1:A:1857:VAL:HG21	1:A:1913:THR:OG1	2.13	0.49
1:A:1991:ILE:HD13	1:A:2008:LEU:HD11	1.94	0.49
2:B:74:LYS:NZ	4:B:403:HOH:O	2.46	0.48
2:B:228:PHE:CD2	2:B:271:GLN:HG2	2.49	0.48
1:A:1859:ARG:NH1	1:A:1979[B]:MET:HE3	2.29	0.48
2:B:54[A]:MET:HE2	4:B:442:HOH:O	2.13	0.48
1:A:1844:PHE:O	1:A:1885:LYS:NZ	2.36	0.47
2:B:54[A]:MET:HE1	4:B:438:HOH:O	2.14	0.47
1:A:1962:ARG:HD3	1:A:1962:ARG:H	1.80	0.46
1:A:1991:ILE:HG21	1:A:2008:LEU:HD22	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1942:ASP:HB2	1:A:1943:PRO:HD3	1.98	0.45
2:B:187:GLU:H	2:B:187:GLU:CD	2.25	0.45
2:B:258:LYS:HD2	2:B:258:LYS:H	1.82	0.45
1:A:2060:LEU:O	1:A:2063:TYR:HB3	2.17	0.44
1:A:2058:LEU:C	1:A:2058:LEU:HD23	2.43	0.44
2:B:43:VAL:O	2:B:43:VAL:CG1	2.65	0.44
2:B:17:ILE:O	2:B:17:ILE:HG23	2.18	0.44
2:B:1:MET:HE3	2:B:35:ASP:O	2.17	0.43
2:B:190:ARG:HG3	2:B:203[B]:TYR:CZ	2.53	0.43
1:A:1974:LEU:HD23	1:A:1989:PHE:CZ	2.51	0.43
2:B:70:GLN:HB3	2:B:81:MET:HE2	1.98	0.43
1:A:1857:VAL:HG13	1:A:1894:ILE:HG13	2.00	0.43
1:A:1992:TYR:O	1:A:1993:ASP:C	2.60	0.43
2:B:85:ASP:C	2:B:85:ASP:OD1	2.62	0.43
2:B:61:ASP:OD1	2:B:61:ASP:C	2.61	0.43
2:B:69:ILE:HD13	2:B:80:MET:HA	2.01	0.42
2:B:77:LEU:N	2:B:77:LEU:HD23	2.34	0.42
2:B:260:MET:HE2	2:B:260:MET:HB3	1.88	0.42
1:A:1854:ASP:CG	1:A:1937:ARG:HH21	2.28	0.42
1:A:1991:ILE:O	1:A:1991:ILE:HG23	2.20	0.42
1:A:2028:SER:O	1:A:2030:PRO:HD3	2.19	0.42
1:A:1835:MET:H	1:A:1959:THR:HA	1.85	0.41
2:B:68:TYR:CE2	2:B:81:MET:HE3	2.55	0.41
2:B:77:LEU:HD23	2:B:77:LEU:H	1.85	0.41
2:B:96:PHE:CB	2:B:102:MET:HE3	2.52	0.40
1:A:1944:LEU:HD12	1:A:1944:LEU:HA	1.80	0.40
2:B:190:ARG:HG3	2:B:203[B]:TYR:CE2	2.56	0.40
2:B:214:PHE:O	2:B:215:LYS:HB2	2.21	0.40
2:B:77:LEU:N	2:B:77:LEU:CD2	2.84	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	
1	A	258/258 (100%)	248 (96%)	9 (4%)	1 (0%)	30	16
2	B	315/308 (102%)	305 (97%)	9 (3%)	1 (0%)	36	24
All	All	573/566 (101%)	553 (96%)	18 (3%)	2 (0%)	30	24

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2061	THR
2	B	106	PRO

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	237/233 (102%)	233 (98%)	4 (2%)	53	31
2	B	294/284 (104%)	290 (99%)	4 (1%)	59	40
All	All	531/517 (103%)	523 (98%)	8 (2%)	59	36

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1835	MET
1	A	1962	ARG
1	A	1991	ILE
1	A	2066	LYS
2	B	45[A]	HIS
2	B	45[B]	HIS
2	B	77	LEU
2	B	281	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1863	HIS

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Mol	Chain	Res	Type
1	A	1869	ASN
1	A	2038	HIS
2	B	47	GLN
2	B	259	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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	VGU	A	2101	-	12,14,14	4.54	5 (41%)	15,17,17	2.00	7 (46%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	VGU	A	2101	-	-	4/9/11/11	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	2101	VGU	C7-N	13.87	1.46	1.29
3	A	2101	VGU	C4-C3	-4.59	1.32	1.39
3	A	2101	VGU	C1-C2	-3.16	1.33	1.38
3	A	2101	VGU	O2-N1	-2.59	1.33	1.40
3	A	2101	VGU	C2-C3	-2.57	1.35	1.39

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	2101	VGU	C3-C6-N	3.47	120.13	114.89
3	A	2101	VGU	C7-N-C6	-3.42	116.19	121.14
3	A	2101	VGU	C2-C3-C6	-2.60	115.77	120.45
3	A	2101	VGU	C8-C7-N	-2.58	116.56	122.52
3	A	2101	VGU	C2-C3-C4	2.49	121.73	118.57
3	A	2101	VGU	C1-C2-C3	-2.13	118.27	120.36
3	A	2101	VGU	O-C6-N	-2.09	120.05	124.73

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	2101	VGU	C7-C8-N1-O2
3	A	2101	VGU	O1-C8-N1-O2
3	A	2101	VGU	N-C7-C8-N1
3	A	2101	VGU	N-C7-C8-O1

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 [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	A	236/258 (91%)	2.19	109 (46%) 0 1	26, 72, 128, 198	11 (4%)
2	B	300/308 (97%)	2.19	142 (47%) 0 1	23, 75, 137, 230	9 (3%)
All	All	536/566 (94%)	2.19	251 (46%) 0 1	23, 74, 131, 230	20 (3%)

All (251) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1878	CYS	8.4
2	B	203[A]	TYR	8.2
1	A	1989	PHE	7.9
1	A	2069	VAL	7.5
1	A	2060	LEU	6.4
2	B	152	LEU	6.2
2	B	246	MET	5.9
1	A	2040	TRP	5.8
1	A	2034	ILE	5.6
1	A	1988	LEU	5.6
2	B	256	VAL	5.5
1	A	1951	PHE	5.3
2	B	181	ILE	5.3
2	B	283	LEU	5.3
1	A	1940	MET	5.2
1	A	2015	LEU	5.2
1	A	2039	LEU	5.2
2	B	293	LEU	5.1
2	B	54[A]	MET	5.1
2	B	226	PHE	5.0
2	B	288	VAL	5.0
2	B	101	MET	5.0
2	B	316	LEU	5.0
2	B	170	SER	5.0

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Mol	Chain	Res	Type	RSRZ
1	A	2037	TYR	4.9
1	A	2036	SER	4.8
1	A	2058	LEU	4.8
1	A	2068	ASN	4.7
2	B	172	PRO	4.7
2	B	183	PHE	4.7
1	A	1898[A]	VAL	4.6
2	B	245	ALA	4.5
2	B	108	ILE	4.5
2	B	111	ASP	4.4
2	B	126	ILE	4.4
2	B	214	PHE	4.4
1	A	2059	ILE	4.4
2	B	92	ILE	4.4
2	B	1	MET	4.4
1	A	1963	LEU	4.3
1	A	2026	LEU	4.3
2	B	275	LEU	4.3
1	A	2043	PHE	4.3
1	A	2061	THR	4.2
1	A	2067	TYR	4.2
2	B	237	TYR	4.2
2	B	38	ILE	4.2
2	B	213	ILE	4.2
2	B	308	ILE	4.2
1	A	2063	TYR	4.1
2	B	294	TYR	4.1
2	B	227	ALA	4.0
1	A	1833	GLY	4.0
2	B	313	TYR	4.0
1	A	2017[A]	THR	3.9
1	A	1969	MET	3.9
1	A	1958	PRO	3.9
2	B	317	LEU	3.8
1	A	1841	ALA	3.8
2	B	284	LEU	3.7
1	A	1834	ALA	3.7
2	B	122[A]	GLN	3.7
2	B	206	ASN	3.7
1	A	2051	ILE	3.6
2	B	276	PRO	3.6
1	A	2004	ALA	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	2038	HIS	3.6
1	A	2048	TRP	3.5
2	B	73	PRO	3.5
2	B	118	THR	3.5
1	A	2064	GLY	3.5
2	B	299	LYS	3.5
1	A	1860	VAL	3.5
2	B	53	SER	3.5
1	A	2033	THR	3.5
1	A	2049	ILE	3.5
2	B	17	ILE	3.4
1	A	1965	PHE	3.4
1	A	2065	ARG	3.4
2	B	193	HIS	3.4
1	A	2027	LEU	3.4
2	B	137	PHE	3.4
2	B	174	HIS	3.3
2	B	291	ILE	3.3
2	B	210	LEU	3.3
1	A	1927	GLU	3.3
1	A	1840	TYR	3.3
1	A	2025	ILE	3.3
1	A	2028	SER	3.3
2	B	204[A]	TYR	3.2
2	B	255	THR	3.2
1	A	1995	TRP	3.2
1	A	2012	LEU	3.2
2	B	173	ALA	3.2
2	B	93	VAL	3.2
2	B	106	PRO	3.1
2	B	261	LEU	3.1
2	B	134	GLU	3.1
2	B	289	TRP	3.1
1	A	1875	ILE	3.1
2	B	282	ILE	3.1
2	B	208	VAL	3.1
2	B	77	LEU	3.1
2	B	207	THR	3.1
2	B	30	PHE	3.1
2	B	314	PRO	3.0
2	B	71	TYR	3.0
2	B	24	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
2	B	279	TYR	3.0
2	B	87	ALA	3.0
2	B	10	PRO	3.0
1	A	1977	VAL	3.0
1	A	1835	MET	3.0
1	A	2030	PRO	3.0
1	A	1866	PHE	3.0
1	A	1837	SER	2.9
1	A	1970	SER	2.9
2	B	244	HIS	2.9
2	B	221	PHE	2.9
2	B	217	SER	2.9
2	B	146	THR	2.9
1	A	2010	LEU	2.9
2	B	128	LYS	2.9
1	A	1992	TYR	2.9
2	B	224	LEU	2.9
2	B	40	HIS	2.9
2	B	45[A]	HIS	2.9
2	B	254	ALA	2.8
1	A	1961[A]	LEU	2.8
2	B	258	LYS	2.8
2	B	280	SER	2.8
2	B	29	PRO	2.8
2	B	220	TYR	2.8
1	A	1873	LYS	2.8
2	B	21	SER	2.8
2	B	234	PHE	2.8
2	B	69	ILE	2.8
2	B	96	PHE	2.7
2	B	64[A]	MET	2.7
1	A	1874	ALA	2.7
2	B	22	PHE	2.7
2	B	198	PHE	2.7
1	A	1862	VAL	2.7
1	A	1996	LEU	2.7
1	A	1971	ILE	2.7
1	A	2035	LYS	2.7
2	B	178[A]	TYR	2.7
1	A	1980[A]	LYS	2.7
1	A	2055	MET	2.7
2	B	50	ASP	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	1890	PHE	2.6
1	A	2050	THR	2.6
2	B	121	VAL	2.6
2	B	114	TRP	2.6
2	B	277	GLU	2.6
1	A	2044	THR	2.6
2	B	14	THR	2.6
1	A	1924	LEU	2.6
1	A	1998	ARG	2.6
2	B	171	ASP	2.6
2	B	78	TYR	2.6
1	A	2041	PRO	2.6
2	B	145	THR	2.6
2	B	281	ASP	2.6
1	A	1848	ILE	2.6
1	A	2001[A]	SER	2.5
2	B	199	LEU	2.5
2	B	180	VAL	2.5
1	A	2042	SER	2.5
1	A	1967	ALA	2.5
2	B	131	ARG	2.5
2	B	292	CYS	2.5
2	B	20	TYR	2.5
1	A	1863	HIS	2.5
1	A	1987	VAL	2.5
2	B	267	ILE	2.5
2	B	99	ARG	2.5
1	A	1905	LEU	2.4
1	A	1947	HIS	2.4
1	A	1935	VAL	2.4
1	A	1999	ILE	2.4
2	B	129	ILE	2.4
1	A	1899[A]	TRP	2.4
2	B	285	ASN	2.4
2	B	191	PRO	2.4
1	A	1939	ALA	2.4
2	B	82	GLU	2.4
1	A	1956	ILE	2.4
1	A	1917	VAL	2.4
2	B	48	HIS	2.4
2	B	306	GLU	2.3
1	A	1968	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
2	B	49	ALA	2.3
2	B	287	ARG	2.3
2	B	46	PHE	2.3
2	B	243	TRP	2.3
2	B	28	GLN	2.3
2	B	43	VAL	2.3
1	A	2066	LYS	2.3
1	A	1936	THR	2.3
1	A	1920	LEU	2.3
2	B	89	PHE	2.3
2	B	123[A]	MET	2.3
2	B	259	HIS	2.3
1	A	1891	LEU	2.3
1	A	1846	ASN	2.3
1	A	1838	SER	2.3
1	A	2023	LYS	2.3
2	B	307	LYS	2.3
1	A	1950	ASP	2.2
1	A	2032	ILE	2.2
2	B	94	HIS	2.2
2	B	13	VAL	2.2
2	B	286	GLU	2.2
1	A	1836	ASN	2.2
1	A	1910	LYS	2.2
1	A	1941	LEU	2.2
1	A	1865	THR	2.2
2	B	2	ASN	2.2
2	B	27	ASN	2.2
2	B	139	TYR	2.2
2	B	211[A]	GLN	2.1
2	B	301	SER	2.1
1	A	2003	THR	2.1
2	B	6	PHE	2.1
2	B	233	PHE	2.1
1	A	1893	ILE	2.1
2	B	56	TYR	2.1
2	B	219	ASN	2.1
2	B	59	TRP	2.1
2	B	151	GLU	2.1
2	B	278	GLN	2.1
2	B	228	PHE	2.1
1	A	1933	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
2	B	270	TYR	2.1
2	B	272	ILE	2.1
1	A	1901	GLY	2.1
1	A	1870	VAL	2.1
2	B	103	VAL	2.1
2	B	218	SER	2.1
1	A	1962	ARG	2.1
2	B	127	ARG	2.1
1	A	1964	PRO	2.1
1	A	1931[A]	LYS	2.1
2	B	132	LYS	2.1
1	A	1851	PHE	2.1
2	B	52	SER	2.0
1	A	1888	HIS	2.0
1	A	1864	LYS	2.0
2	B	107	LYS	2.0
2	B	147	VAL	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	VGU	A	2101	14/14	0.63	0.18	20,20,20,20	14

6.5 Other polymers [i](#)

There are no such residues in this entry.