



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 27, 2026 – 05:20 PM EDT

PDB ID : 7FM2 / pdb_00007fm2
Title : PanDDA analysis group deposition – Aar2/RNaseH in complex with fragment P05H04 from the F2X-Universal Library
Authors : Barthel, T.; Wollenhaupt, J.; Lima, G.M.A.; Wahl, M.C.; Weiss, M.S.
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

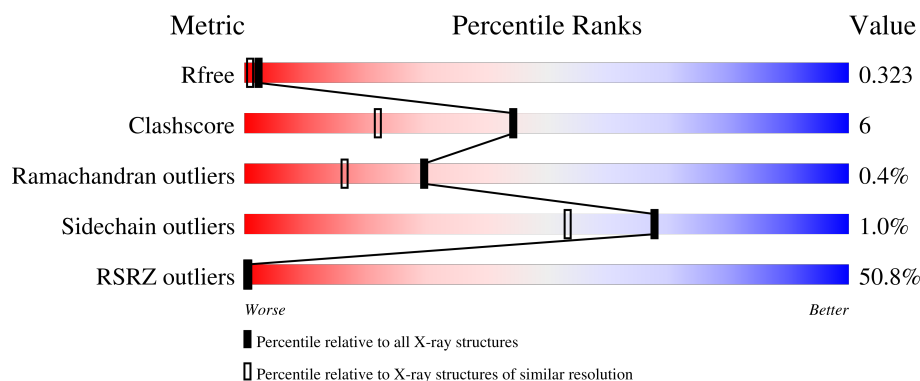
1 Overall quality at a glance

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	
2	B	308	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9189 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	0	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:

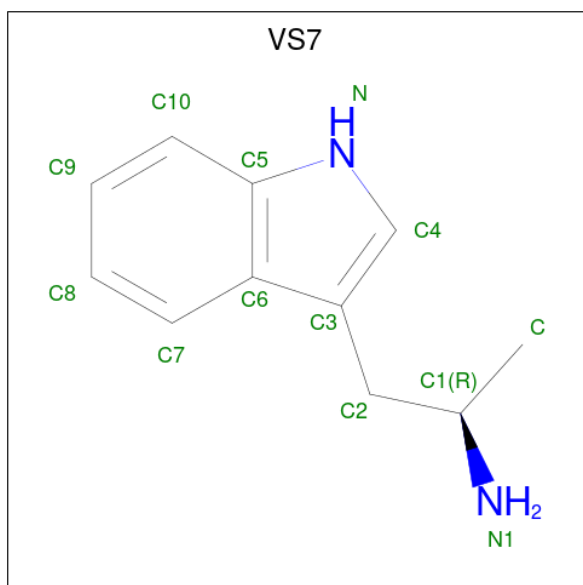
Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	GLY	-	expression tag	UNP P32357
B	-2	ALA	-	expression tag	UNP P32357
B	-1	MET	-	expression tag	UNP P32357
B	0	ALA	-	expression tag	UNP P32357
B	166	SER	LEU	conflict	UNP P32357
B	167	SER	LYS	conflict	UNP P32357
B	?	-	LEU	deletion	UNP P32357
B	?	-	GLN	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 (2R)-1-(1H-indol-3-yl)propan-2-amine (CCD ID: VS7) (formula: C₁₁H₁₄N₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	N	0	0
			13	11	2		

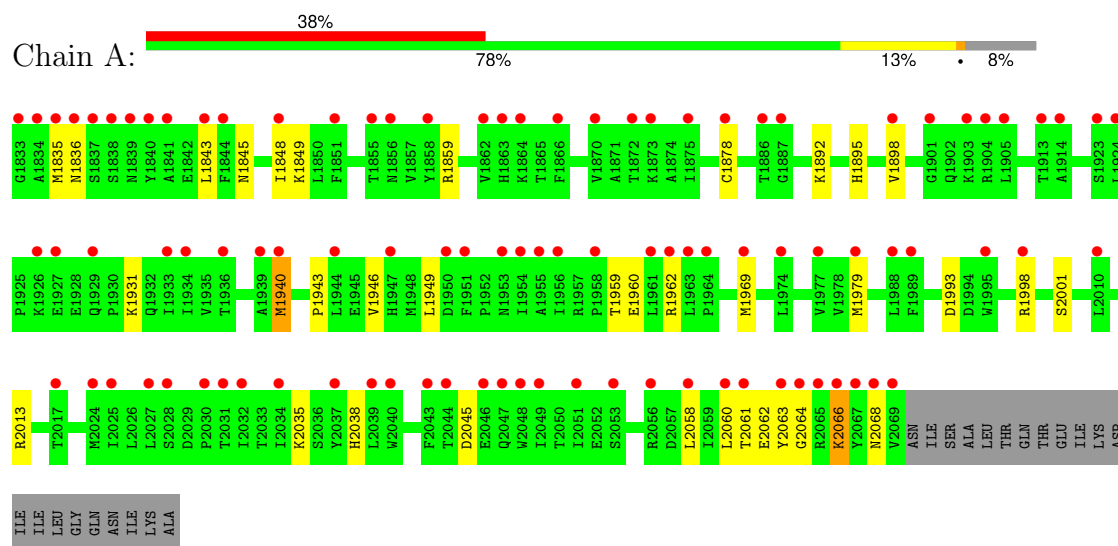
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	32	Total	O	0	0
			32	32		
4	B	32	Total	O	0	0
			32	32		

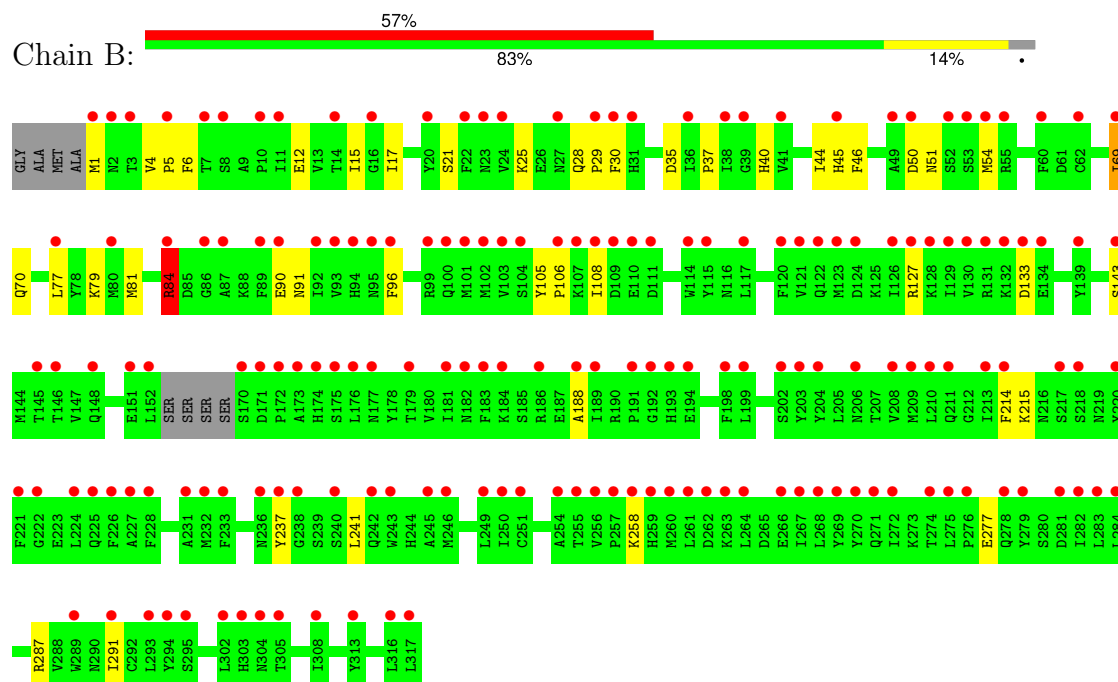
3 Residue-property plots

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	88.01Å 81.49Å 92.93Å 90.00° 107.80° 90.00°	Depositor
Resolution (Å)	44.24 – 1.72 44.24 – 1.72	Depositor EDS
% Data completeness (in resolution range)	99.5 (44.24-1.72) 99.9 (44.24-1.72)	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}	0.275 , 0.319 0.283 , 0.323	Depositor DCC
R_{free} test set	2101 reflections (3.16%)	wwPDB-VP
Wilson B-factor (Å ²)	40.2	Xtriage
Anisotropy	0.410	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 48.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9189	wwPDB-VP
Average B, all atoms (Å ²)	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.12% 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: VS7

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	0.71	1/2149 (0.0%)	0.81	0/2911
2	B	0.80	2/2739 (0.1%)	0.95	11/3699 (0.3%)
All	All	0.76	3/4888 (0.1%)	0.89	11/6610 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	5	PRO	N-CA	16.49	1.67	1.47
2	B	4	VAL	C-N	8.72	1.45	1.33
1	A	1940	MET	SD-CE	-5.90	1.64	1.79

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	4	VAL	CA-C-N	14.53	134.73	119.76
2	B	4	VAL	C-N-CA	14.53	134.73	119.76
2	B	5	PRO	CA-N-CD	-8.23	100.47	112.00
2	B	5	PRO	N-CA-C	-8.11	98.49	111.14
2	B	5	PRO	CB-CA-C	6.24	119.51	111.46
2	B	90	GLU	CB-CA-C	-6.01	101.44	110.88
2	B	96	PHE	CA-CB-CG	5.97	119.77	113.80
2	B	84	ARG	N-CA-C	-5.86	105.96	113.23
2	B	30	PHE	CA-C-N	-5.37	112.73	122.50
2	B	30	PHE	C-N-CA	-5.37	112.73	122.50
2	B	6	PHE	CA-C-O	-5.08	115.32	120.71

There are no chirality outliers.

There are no planarity outliers.

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	A	2008	2060	1974	27	0
2	B	2580	2464	2398	26	0
3	B	13	0	0	1	0
4	A	32	0	0	3	0
4	B	32	0	0	1	0
All	All	4665	4524	4372	54	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 6.

All (54) 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:46:PHE:HZ	2:B:69:ILE:HD12	1.43	0.83
1:A:2063:TYR:O	4:A:2101:HOH:O	1.96	0.81
1:A:2061:THR:O	1:A:2064:GLY:N	2.16	0.78
1:A:2064:GLY:HA2	4:A:2101:HOH:O	1.87	0.74
2:B:46:PHE:CZ	2:B:69:ILE:HD12	2.29	0.64
1:A:1848:ILE:H	1:A:1931[A]:LYS:HZ2	1.46	0.63
1:A:1878:CYS:SG	1:A:1969:MET:HE3	2.41	0.61
2:B:17:ILE:HD13	2:B:44[B]:ILE:CG1	2.34	0.58
2:B:287:ARG:O	2:B:291:ILE:HD13	2.06	0.56
2:B:77:LEU:HD21	2:B:79:LYS:HE3	1.89	0.54
2:B:40:HIS:HD2	4:B:523:HOH:O	1.91	0.53
3:B:401:VS7:C4	3:B:401:VS7:N1	2.72	0.52
1:A:1843:LEU:HA	1:A:1849:LYS:HD2	1.91	0.52
1:A:1962:ARG:O	1:A:2013:ARG:NH1	2.43	0.51
2:B:70:GLN:HB3	2:B:81:MET:HE2	1.92	0.50
1:A:2060:LEU:HB2	4:A:2132:HOH:O	2.11	0.50
2:B:50:ASP:OD1	2:B:51:ASN:N	2.46	0.49
1:A:1845:ASN:OD1	1:A:1849:LYS:NZ	2.46	0.49
1:A:1998:ARG:NH1	1:A:2045:ASP:OD2	2.39	0.49
1:A:1962:ARG:HD3	1:A:1962:ARG:H	1.79	0.48
2:B:84:ARG:HB2	2:B:84:ARG:HH11	1.79	0.47
2:B:277:GLU:CD	2:B:277:GLU:H	2.23	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:258:LYS:HD2	2:B:258:LYS:H	1.80	0.47
1:A:2062:GLU:O	1:A:2066:LYS:HB2	2.16	0.46
1:A:1895:HIS:O	1:A:1898[A]:VAL:HG22	2.16	0.46
1:A:1946:VAL:O	1:A:1949:LEU:HG	2.17	0.45
1:A:1993:ASP:HB3	1:A:2038:HIS:CD2	2.52	0.45
2:B:15:ILE:O	2:B:21:SER:HA	2.17	0.45
1:A:1859:ARG:NH1	1:A:1979[C]:MET:SD	2.90	0.45
1:A:2058:LEU:C	1:A:2058:LEU:HD23	2.41	0.45
2:B:127:ARG:NH1	2:B:133:ASP:O	2.49	0.45
2:B:17:ILE:CD1	2:B:44[B]:ILE:HG13	2.47	0.44
2:B:17:ILE:HD13	2:B:44[B]:ILE:HG13	1.99	0.44
1:A:1940:MET:C	1:A:1943:PRO:HD2	2.42	0.43
2:B:108:ILE:HD13	2:B:108:ILE:N	2.34	0.43
2:B:28:GLN:HG3	2:B:29:PRO:HD2	2.00	0.42
2:B:214:PHE:O	2:B:215:LYS:HB2	2.19	0.42
1:A:1836:ASN:HB3	1:A:1960[A]:GLU:HB2	2.02	0.42
1:A:1878:CYS:HA	1:A:1892:LYS:O	2.19	0.42
1:A:1859:ARG:HH12	1:A:1979[A]:MET:HE3	1.83	0.42
2:B:17:ILE:HD13	2:B:44[B]:ILE:HG12	2.01	0.42
1:A:1835:MET:HE3	1:A:1959:THR:OG1	2.19	0.42
2:B:54[B]:MET:HE1	2:B:143:SER:HB3	2.02	0.41
2:B:1:MET:HB3	2:B:35:ASP:HA	2.03	0.41
1:A:2035:LYS:HD3	1:A:2035:LYS:HA	1.85	0.41
2:B:237:TYR:CE2	2:B:241:LEU:HD11	2.55	0.41
2:B:12:GLU:HG3	2:B:25:LYS:HA	2.03	0.41
1:A:2061:THR:O	1:A:2062:GLU:C	2.64	0.40
2:B:28:GLN:CG	2:B:29:PRO:HD2	2.51	0.40
2:B:37:PRO:HD3	2:B:105:TYR:CD1	2.56	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	258/258 (100%)	248 (96%)	9 (4%)	1 (0%)	30	16
2	B	315/308 (102%)	303 (96%)	11 (4%)	1 (0%)	36	24
All	All	573/566 (101%)	551 (96%)	20 (4%)	2 (0%)	30	24

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2068	ASN
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%)	236 (100%)	1 (0%)	84	76
2	B	294/284 (104%)	289 (98%)	5 (2%)	53	31
All	All	531/517 (103%)	525 (99%)	6 (1%)	68	50

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2066	LYS
2	B	45[A]	HIS
2	B	45[B]	HIS
2	B	69	ILE
2	B	84	ARG
2	B	91	ASN

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

Mol	Chain	Res	Type
1	A	1869	ASN
2	B	2	ASN
2	B	47	GLN

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Mol	Chain	Res	Type
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	VS7	B	401	-	14,14,14	1.18	1 (7%)	19,19,19	0.87	0

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	VS7	B	401	-	-	3/4/4/4	0/2/2/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	401	VS7	C1-N1	2.28	1.55	1.49

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	401	VS7	C-C1-C2-C3
3	B	401	VS7	C1-C2-C3-C6
3	B	401	VS7	C1-C2-C3-C4

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	401	VS7	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 [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	237/258 (91%)	2.05	97 (40%) 0 1	23, 66, 135, 181	12 (5%)
2	B	300/308 (97%)	2.46	176 (58%) 0 0	21, 76, 136, 205	9 (3%)
All	All	537/566 (94%)	2.28	273 (50%) 0 0	21, 72, 136, 205	21 (3%)

All (273) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	203[A]	TYR	19.4
1	A	1979[A]	MET	13.0
1	A	1878	CYS	9.7
1	A	2060	LEU	8.6
2	B	54[A]	MET	8.5
1	A	2069	VAL	7.6
2	B	101	MET	6.8
2	B	316	LEU	6.3
2	B	152	LEU	6.1
2	B	231	ALA	5.7
1	A	2027	LEU	5.4
2	B	282	ILE	5.3
2	B	1	MET	5.2
1	A	1834	ALA	5.1
2	B	170	SER	5.0
2	B	55[A]	ARG	4.9
2	B	183	PHE	4.9
1	A	2067	TYR	4.8
2	B	129	ILE	4.8
1	A	2030	PRO	4.8
2	B	5	PRO	4.8
2	B	275	LEU	4.7
2	B	39	GLY	4.7
2	B	313	TYR	4.6

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Mol	Chain	Res	Type	RSRZ
1	A	2064	GLY	4.6
2	B	226	PHE	4.6
1	A	1862	VAL	4.5
2	B	224	LEU	4.5
2	B	80	MET	4.5
2	B	173	ALA	4.5
2	B	120	PHE	4.5
2	B	176	LEU	4.5
2	B	172	PRO	4.4
2	B	269	TYR	4.4
1	A	1963	LEU	4.3
2	B	111	ASP	4.3
1	A	1955	ALA	4.3
1	A	2063	TYR	4.3
2	B	130	VAL	4.3
2	B	272	ILE	4.3
2	B	206	ASN	4.2
2	B	145	THR	4.2
1	A	1833	GLY	4.1
2	B	267	ILE	4.1
2	B	22	PHE	4.1
2	B	107	LYS	4.1
2	B	289	TRP	4.0
1	A	1974	LEU	4.0
2	B	246	MET	4.0
1	A	2051	ILE	4.0
2	B	213	ILE	4.0
2	B	218	SER	4.0
1	A	2037	TYR	4.0
2	B	237	TYR	4.0
2	B	96	PHE	4.0
2	B	41[A]	VAL	4.0
2	B	208	VAL	3.9
2	B	181	ILE	3.9
1	A	2068	ASN	3.9
2	B	92	ILE	3.8
2	B	294	TYR	3.8
2	B	276	PRO	3.8
1	A	2017[A]	THR	3.8
2	B	30	PHE	3.7
2	B	53	SER	3.7
2	B	262	ASP	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	2065	ARG	3.7
1	A	2039	LEU	3.7
2	B	122[A]	GLN	3.7
2	B	217	SER	3.7
2	B	284	LEU	3.6
1	A	1969	MET	3.6
2	B	126	ILE	3.6
2	B	227	ALA	3.6
2	B	274	THR	3.6
2	B	221	PHE	3.6
2	B	184	LYS	3.6
1	A	2061	THR	3.5
1	A	1988	LEU	3.5
1	A	1872	THR	3.5
2	B	260	MET	3.5
2	B	249	LEU	3.5
2	B	204[A]	TYR	3.5
2	B	214	PHE	3.4
2	B	261	LEU	3.4
2	B	121	VAL	3.4
2	B	256	VAL	3.4
2	B	174	HIS	3.4
2	B	189	ILE	3.4
2	B	283	LEU	3.4
2	B	243	TRP	3.4
1	A	1836	ASN	3.4
1	A	1944	LEU	3.4
1	A	2043	PHE	3.4
2	B	8	SER	3.4
2	B	171	ASP	3.3
2	B	193	HIS	3.3
2	B	31	HIS	3.3
1	A	2024[A]	MET	3.3
1	A	2049	ILE	3.3
2	B	134	GLU	3.3
2	B	291	ILE	3.3
2	B	198	PHE	3.3
1	A	1913	THR	3.2
2	B	308	ILE	3.2
2	B	106	PRO	3.2
2	B	90	GLU	3.2
2	B	194	GLU	3.2

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Mol	Chain	Res	Type	RSRZ
2	B	60	PHE	3.2
2	B	233	PHE	3.2
1	A	2058	LEU	3.2
2	B	50	ASP	3.2
1	A	1840	TYR	3.1
2	B	10	PRO	3.1
2	B	270	TYR	3.1
2	B	279	TYR	3.1
2	B	108	ILE	3.1
1	A	2028	SER	3.1
2	B	87	ALA	3.1
1	A	1956	ILE	3.1
2	B	45[A]	HIS	3.1
2	B	264	LEU	3.1
2	B	317	LEU	3.1
2	B	225	GLN	3.1
1	A	1998	ARG	3.1
2	B	103	VAL	3.1
1	A	1958	PRO	3.1
1	A	2046	GLU	3.1
2	B	11	ILE	3.1
1	A	1841	ALA	3.1
1	A	1837	SER	3.1
2	B	124	ASP	3.0
2	B	255	THR	3.0
2	B	202	SER	3.0
2	B	258	LYS	3.0
1	A	1870	VAL	2.9
1	A	1838	SER	2.9
2	B	7	THR	2.9
2	B	305	THR	2.9
1	A	1848	ILE	2.9
2	B	38	ILE	2.9
1	A	1951	PHE	2.9
1	A	1856[A]	ASN	2.9
1	A	2032	ILE	2.9
1	A	2053[A]	SER	2.9
1	A	1866	PHE	2.9
2	B	268	LEU	2.9
2	B	102	MET	2.9
2	B	94	HIS	2.8
2	B	146	THR	2.8

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Mol	Chain	Res	Type	RSRZ
2	B	303	HIS	2.8
2	B	139	TYR	2.8
2	B	211[A]	GLN	2.8
2	B	186	ARG	2.8
2	B	89	PHE	2.7
1	A	1898[A]	VAL	2.7
1	A	1839	ASN	2.7
1	A	1950	ASP	2.7
2	B	240	SER	2.7
2	B	123[A]	MET	2.7
1	A	1863	HIS	2.7
1	A	2044	THR	2.7
1	A	1835	MET	2.7
2	B	191	PRO	2.6
2	B	52	SER	2.6
2	B	131	ARG	2.6
1	A	1855[A]	THR	2.6
1	A	1989	PHE	2.6
1	A	2048	TRP	2.6
1	A	2031	THR	2.6
1	A	1929	GLN	2.6
1	A	1873	LYS	2.6
1	A	1875	ILE	2.6
2	B	209	MET	2.5
1	A	1953	ASN	2.5
2	B	2	ASN	2.5
2	B	100	GLN	2.5
2	B	251	CYS	2.5
1	A	1843	LEU	2.5
2	B	257	PRO	2.5
2	B	245	ALA	2.5
1	A	1933	ILE	2.5
1	A	2034	ILE	2.5
1	A	1927	GLU	2.5
2	B	236	ASN	2.4
1	A	2040	TRP	2.4
1	A	1905	LEU	2.4
2	B	16	GLY	2.4
1	A	2047	GLN	2.4
2	B	109	ASP	2.4
1	A	1961[A]	LEU	2.4
1	A	2010	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	29	PRO	2.4
2	B	77	LEU	2.4
2	B	210	LEU	2.4
2	B	49	ALA	2.4
2	B	175	SER	2.4
2	B	254	ALA	2.4
2	B	143	SER	2.4
2	B	302	LEU	2.4
1	A	1858	TYR	2.4
2	B	220	TYR	2.4
2	B	232	MET	2.4
2	B	23	ASN	2.3
2	B	188	ALA	2.3
2	B	132	LYS	2.3
2	B	69	ILE	2.3
1	A	1923	SER	2.3
1	A	2066	LYS	2.3
1	A	1940	MET	2.3
1	A	1924	LEU	2.3
1	A	2025	ILE	2.3
2	B	114	TRP	2.3
2	B	3	THR	2.3
2	B	62	CYS	2.3
2	B	104	SER	2.3
2	B	24	VAL	2.3
2	B	14	THR	2.2
1	A	1844	PHE	2.2
1	A	1887	GLY	2.2
2	B	228	PHE	2.2
2	B	115	TYR	2.2
2	B	250	ILE	2.2
2	B	127	ARG	2.2
2	B	304	ASN	2.2
1	A	1939	ALA	2.2
1	A	1901	GLY	2.2
2	B	271	GLN	2.2
2	B	182	ASN	2.2
2	B	263	LYS	2.2
2	B	133	ASP	2.2
1	A	1995	TRP	2.2
2	B	278	GLN	2.2
2	B	110	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
2	B	192	GLY	2.2
1	A	1962	ARG	2.2
1	A	1851	PHE	2.2
1	A	1926	LYS	2.2
1	A	1904	ARG	2.1
2	B	238	GLY	2.1
2	B	199	LEU	2.1
2	B	281	ASP	2.1
2	B	151	GLU	2.1
1	A	1977	VAL	2.1
2	B	27	ASN	2.1
2	B	295	SER	2.1
1	A	1936	THR	2.1
1	A	1934	ILE	2.1
1	A	1954	ILE	2.1
2	B	259	HIS	2.1
2	B	266	GLU	2.1
2	B	293	LEU	2.1
2	B	179	THR	2.1
1	A	1964	PRO	2.1
1	A	1947	HIS	2.1
2	B	36	ILE	2.1
2	B	222	GLY	2.0
2	B	148	GLN	2.0
1	A	2056[A]	ARG	2.0
2	B	93	VAL	2.0
2	B	99	ARG	2.0
2	B	117	LEU	2.0
1	A	1886	THR	2.0
1	A	1864	LYS	2.0
1	A	1903	LYS	2.0
2	B	128	LYS	2.0
2	B	86	GLY	2.0
2	B	95	ASN	2.0
2	B	177[A]	ASN	2.0
2	B	242	GLN	2.0
2	B	20	TYR	2.0
2	B	84	ARG	2.0
1	A	1914	ALA	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	VS7	B	401	13/13	0.78	0.14	20,20,20,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.