



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 18, 2021 – 05:02 am BST

PDB ID : 7A1V
Title : Crystal structure of YTHDF2 in complex with m1A
Authors : Wang, X.; Caffisch, A.
Deposited on : 2020-08-14
Resolution : 2.20 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.23.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.23.1

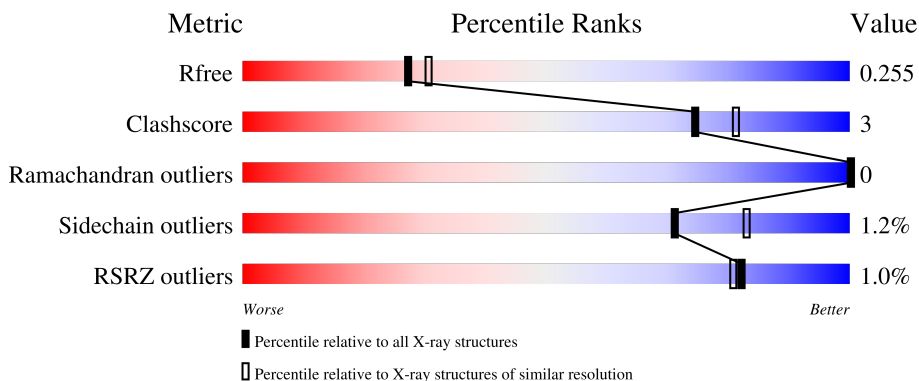
1 Overall quality at a glance

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}	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	167	 2% 83% 6% 11%
1	B	167	 79% 10% 11%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2492 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 YTH domain-containing family protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	149	1189	753	214	217	5	0	0	0
1	B	149	1184	747	211	220	6	0	1	0

There are 44 discrepancies between the modelled and reference sequences:

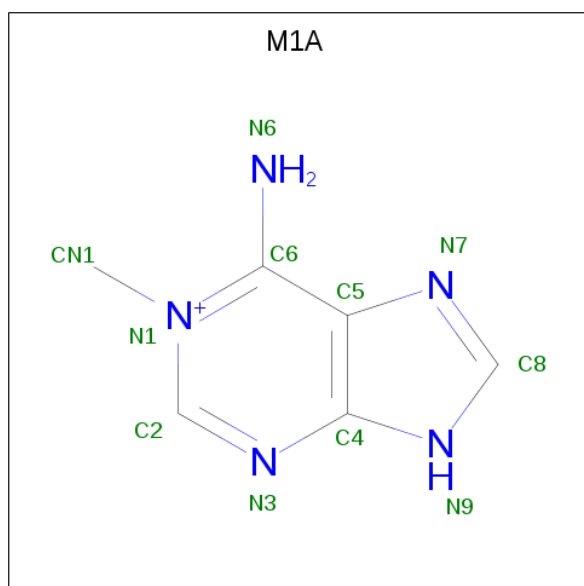
Chain	Residue	Modelled	Actual	Comment	Reference
A	386	MET	-	initiating methionine	UNP Q9Y5A9
A	387	GLY	-	expression tag	UNP Q9Y5A9
A	388	SER	-	expression tag	UNP Q9Y5A9
A	389	SER	-	expression tag	UNP Q9Y5A9
A	390	TYR	-	expression tag	UNP Q9Y5A9
A	391	HIS	-	expression tag	UNP Q9Y5A9
A	392	HIS	-	expression tag	UNP Q9Y5A9
A	393	HIS	-	expression tag	UNP Q9Y5A9
A	394	HIS	-	expression tag	UNP Q9Y5A9
A	395	HIS	-	expression tag	UNP Q9Y5A9
A	396	HIS	-	expression tag	UNP Q9Y5A9
A	397	SER	-	expression tag	UNP Q9Y5A9
A	398	SER	-	expression tag	UNP Q9Y5A9
A	399	GLY	-	expression tag	UNP Q9Y5A9
A	400	GLU	-	expression tag	UNP Q9Y5A9
A	401	ASN	-	expression tag	UNP Q9Y5A9
A	402	LEU	-	expression tag	UNP Q9Y5A9
A	403	TYR	-	expression tag	UNP Q9Y5A9
A	404	PHE	-	expression tag	UNP Q9Y5A9
A	405	GLN	-	expression tag	UNP Q9Y5A9
A	406	HIS	-	expression tag	UNP Q9Y5A9
A	407	MET	-	expression tag	UNP Q9Y5A9
B	386	MET	-	initiating methionine	UNP Q9Y5A9
B	387	GLY	-	expression tag	UNP Q9Y5A9
B	388	SER	-	expression tag	UNP Q9Y5A9

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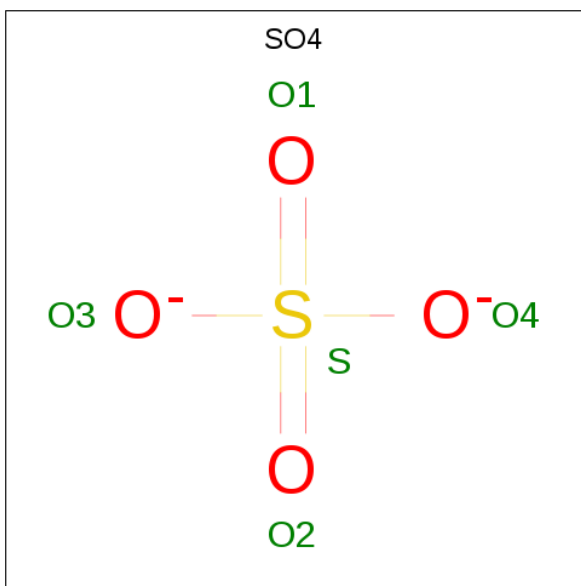
Chain	Residue	Modelled	Actual	Comment	Reference
B	389	SER	-	expression tag	UNP Q9Y5A9
B	390	TYR	-	expression tag	UNP Q9Y5A9
B	391	HIS	-	expression tag	UNP Q9Y5A9
B	392	HIS	-	expression tag	UNP Q9Y5A9
B	393	HIS	-	expression tag	UNP Q9Y5A9
B	394	HIS	-	expression tag	UNP Q9Y5A9
B	395	HIS	-	expression tag	UNP Q9Y5A9
B	396	HIS	-	expression tag	UNP Q9Y5A9
B	397	SER	-	expression tag	UNP Q9Y5A9
B	398	SER	-	expression tag	UNP Q9Y5A9
B	399	GLY	-	expression tag	UNP Q9Y5A9
B	400	GLU	-	expression tag	UNP Q9Y5A9
B	401	ASN	-	expression tag	UNP Q9Y5A9
B	402	LEU	-	expression tag	UNP Q9Y5A9
B	403	TYR	-	expression tag	UNP Q9Y5A9
B	404	PHE	-	expression tag	UNP Q9Y5A9
B	405	GLN	-	expression tag	UNP Q9Y5A9
B	406	HIS	-	expression tag	UNP Q9Y5A9
B	407	MET	-	expression tag	UNP Q9Y5A9

- Molecule 2 is 6-AMINO-1-METHYLPURINE (three-letter code: M1A) (formula: C₆H₈N₅) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	0	0
			11	6	5		

- Molecule 3 is SULFATE ION (three-letter code: 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	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		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	C O	0	0
			6	3 3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	43	Total	O	0	0
			43	43		
5	B	24	Total	O	0	0
			24	24		

4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	79.78Å 79.78Å 113.56Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.87 – 2.20 43.87 – 2.20	Depositor EDS
% Data completeness (in resolution range)	100.0 (43.87-2.20) 97.5 (43.87-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.30 (at 2.20Å)	Xtrriage
Refinement program	PHENIX 1.18.2_3874+SVN	Depositor
R, R_{free}	0.222 , 0.254 0.222 , 0.255	Depositor DCC
R_{free} test set	1040 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	47.9	Xtrriage
Anisotropy	0.033	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 49.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.063 for h,-h-k,-l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2492	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.30% 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: SO4, M1A, GOL

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	0.25	0/1219	0.43	0/1650
1	B	0.25	0/1214	0.43	0/1648
All	All	0.25	0/2433	0.43	0/3298

There are no bond length outliers.

There are no bond angle outliers.

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	1189	0	1128	6	0
1	B	1184	0	1094	9	0
2	A	11	0	8	1	0
3	A	20	0	0	0	0
3	B	15	0	0	0	0
4	B	6	0	8	1	0
5	A	43	0	0	0	0
5	B	24	0	0	0	0
All	All	2492	0	2238	15	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 3.

The worst 5 of 15 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:433:CYS:O	2:A:701:M1A:H2	1.86	0.75
1:A:503:LYS:NZ	1:A:547:TYR:O	2.33	0.60
1:B:426:SER:HG	1:B:499:TRP:HE1	1.52	0.58
1:B:491:TRP:HE1	4:B:701:GOL:H11	1.69	0.58
1:B:475:SER:HB3	1:B:496:ASP:HB2	1.89	0.55

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	147/167 (88%)	143 (97%)	4 (3%)	0	100	100
1	B	148/167 (89%)	144 (97%)	4 (3%)	0	100	100
All	All	295/334 (88%)	287 (97%)	8 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	125/149 (84%)	124 (99%)	1 (1%)	81	90

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	123/149 (83%)	121 (98%)	2 (2%)	62	76
All	All	248/298 (83%)	245 (99%)	3 (1%)	71	83

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	475	SER
1	B	407	MET
1	B	475	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	409	HIS
1	B	539	GLN

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 monosaccharides 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	705	-	4,4,4	0.52	0	6,6,6	0.09	0
3	SO4	A	703	-	4,4,4	0.14	0	6,6,6	0.05	0
3	SO4	A	702	-	4,4,4	0.14	0	6,6,6	0.05	0
4	GOL	B	701	-	5,5,5	0.87	0	5,5,5	1.03	0
2	M1A	A	701	-	9,12,12	1.33	1 (11%)	3,17,17	2.00	2 (66%)
3	SO4	A	704	-	4,4,4	0.14	0	6,6,6	0.04	0
3	SO4	B	704	-	4,4,4	0.52	0	6,6,6	0.06	0
3	SO4	B	702	-	4,4,4	0.14	0	6,6,6	0.05	0
3	SO4	B	703	-	4,4,4	0.14	0	6,6,6	0.05	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
2	M1A	A	701	-	-	-	0/2/2/2
4	GOL	B	701	-	-	4/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	701	M1A	C6-N6	2.49	1.45	1.35

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	701	M1A	C2-N3-C4	2.63	119.88	116.58
2	A	701	M1A	C4-C5-N7	-2.25	107.05	109.40

There are no chirality outliers.

All (4) torsion outliers are listed below:

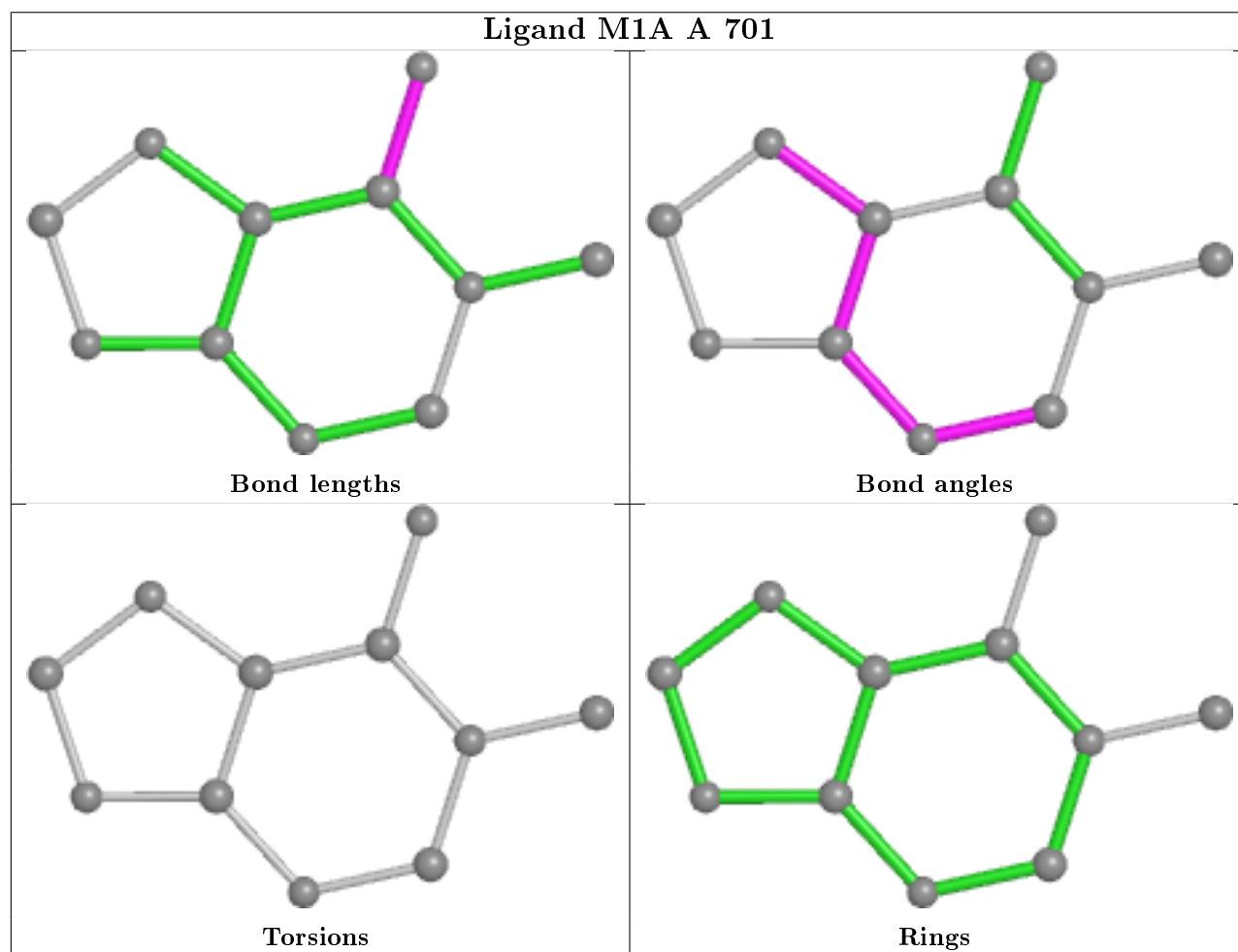
Mol	Chain	Res	Type	Atoms
4	B	701	GOL	C1-C2-C3-O3
4	B	701	GOL	O2-C2-C3-O3
4	B	701	GOL	O1-C1-C2-O2
4	B	701	GOL	O1-C1-C2-C3

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	701	GOL	1	0
2	A	701	M1A	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers

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	149/167 (89%)	-0.04	3 (2%) 65 63	39, 51, 82, 109	0
1	B	149/167 (89%)	0.03	0 100 100	36, 57, 84, 105	0
All	All	298/334 (89%)	-0.01	3 (1%) 82 81	36, 53, 84, 109	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	402	LEU	3.1
1	A	418	TYR	2.1
1	A	492	LYS	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 monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GOL	B	701	6/6	0.64	0.23	71,75,79,81	0
3	SO4	A	703	5/5	0.80	0.16	125,126,129,130	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	SO4	B	703	5/5	0.84	0.17	126,127,129,130	0
3	SO4	B	704	5/5	0.87	0.16	67,73,75,97	0
3	SO4	A	704	5/5	0.92	0.27	92,96,100,109	0
2	M1A	A	701	11/11	0.93	0.16	41,45,48,48	0
3	SO4	A	702	5/5	0.94	0.15	96,97,106,109	0
3	SO4	A	705	5/5	0.95	0.13	56,63,66,89	0
3	SO4	B	702	5/5	0.96	0.10	75,84,87,91	0

6.5 Other polymers [i](#)

There are no such residues in this entry.