



# Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 12:08 pm BST

PDB ID : 6HH5  
Title : ADP-ribosylserine hydrolase ARH3 of *Latimeria chalumnae* in complex with ADP-HPM  
Authors : Ariza, A.  
Deposited on : 2018-08-24  
Resolution : 1.95 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](mailto: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	:	2.11
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)	:	2.11

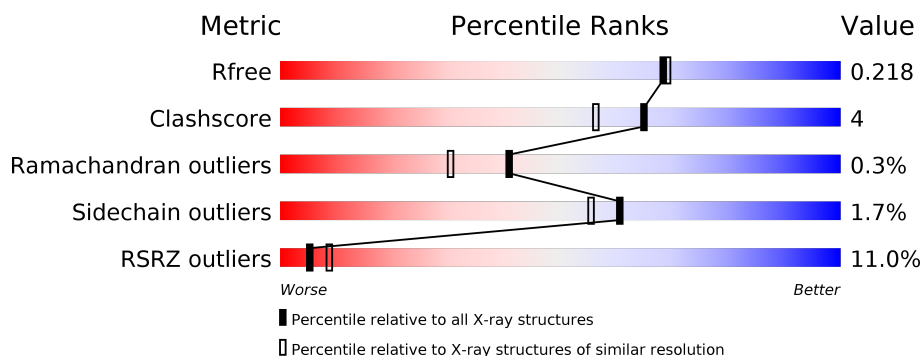
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.95 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	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	350	<div> <div>15%</div> <div>82%</div> <div>10%</div> <div>7%</div> </div>
1	B	350	<div> <div>6%</div> <div>85%</div> <div>9%</div> <div>6%</div> </div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5446 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 ADP-ribosylhydrolase like 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	327	Total	C	N	O	S	0	6	0
			2546	1620	418	493	15			
1	B	328	Total	C	N	O	S	0	7	0
			2561	1631	420	495	15			

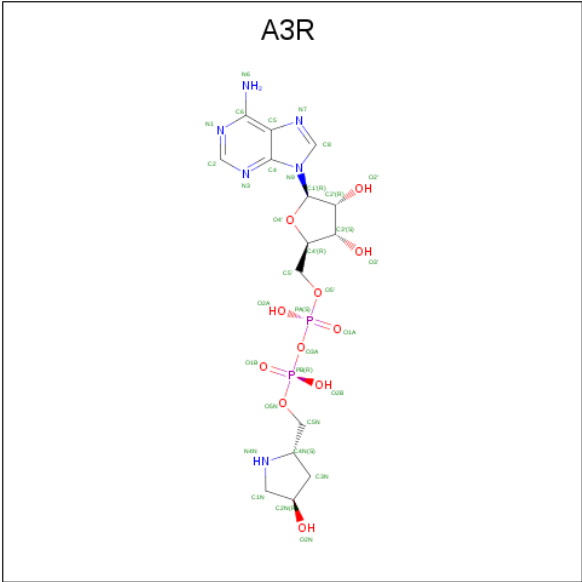
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	7	GLY	-	expression tag	UNP H3BCW1
A	8	PRO	-	expression tag	UNP H3BCW1
A	9	MET	-	expression tag	UNP H3BCW1
B	7	GLY	-	expression tag	UNP H3BCW1
B	8	PRO	-	expression tag	UNP H3BCW1
B	9	MET	-	expression tag	UNP H3BCW1

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is Adenosine Diphosphate (Hydroxymethyl)pyrrolidine monoalcohol (three-letter code: A3R) (formula: C<sub>15</sub>H<sub>24</sub>N<sub>6</sub>O<sub>11</sub>P<sub>2</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			34	15	6	11	2		
3	B	1	Total	C	N	O	P	0	0
			34	15	6	11	2		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

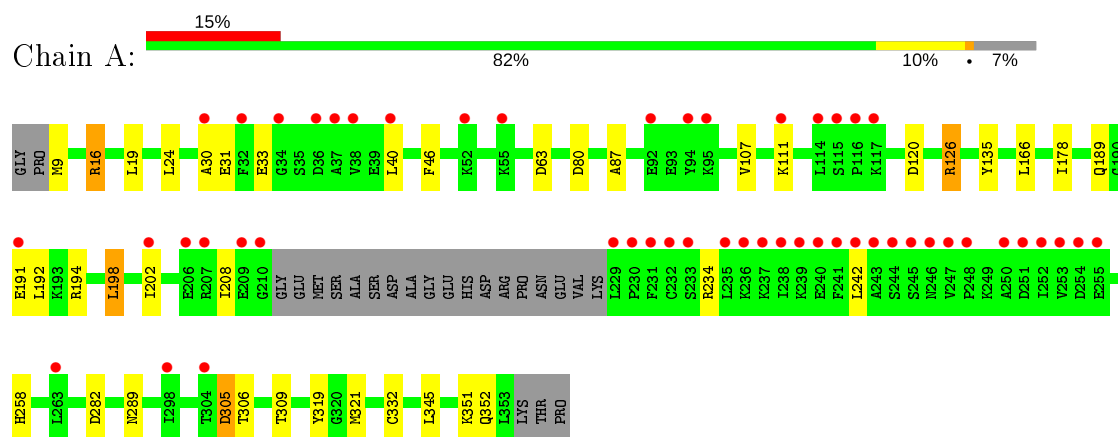
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	86	Total 86	O 86	0	0
5	B	169	Total 171	O 171	0	2

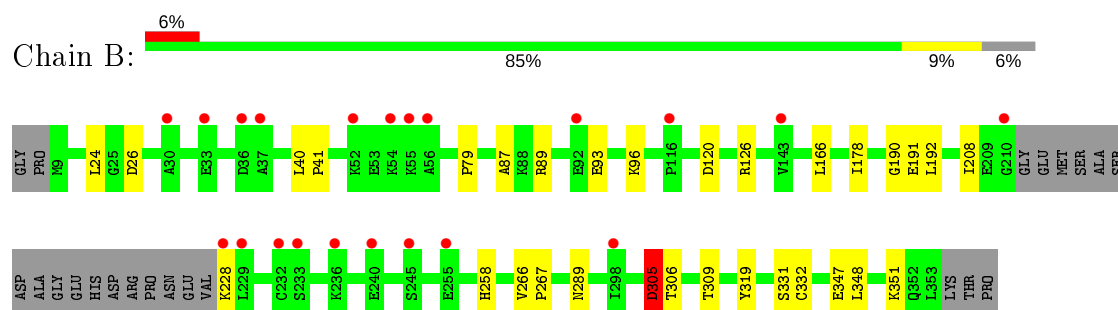
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: ADP-ribosylhydrolase like 2



- Molecule 1: ADP-ribosylhydrolase like 2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	66.87Å 98.18Å 105.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.46 – 1.95 41.43 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.6 (41.46-1.95) 99.6 (41.43-1.95)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.63 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.8.0222	Depositor
R, $R_{free}$	0.186 , 0.214 0.193 , 0.218	Depositor DCC
$R_{free}$ test set	2508 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	38.9	Xtriage
Anisotropy	0.644	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 52.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	5446	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	61.0	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 37.41 % 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 4.3075e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical 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: GOL, MG, A3R

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	0.55	0/2609	0.69	0/3522
1	B	0.61	1/2627 (0.0%)	0.72	2/3544 (0.1%)
All	All	0.58	1/5236 (0.0%)	0.71	2/7066 (0.0%)

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	3
1	B	0	1
All	All	0	4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	93	GLU	CD-OE2	5.24	1.31	1.25

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	305[A]	ASP	N-CA-CB	5.14	119.84	110.60
1	B	305[B]	ASP	N-CA-CB	5.14	119.84	110.60

There are no chirality outliers.

All (4) planarity outliers are listed below:



Mol	Chain	Res	Type	Group
1	A	126	ARG	Sidechain
1	A	16	ARG	Sidechain
1	A	234	ARG	Sidechain
1	B	89	ARG	Sidechain

## 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	2546	0	2543	28	0
1	B	2561	0	2567	14	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	34	0	0	1	0
3	B	34	0	0	0	0
4	A	6	0	8	2	0
4	B	6	0	8	0	0
5	A	86	0	0	3	0
5	B	171	0	0	1	0
All	All	5446	0	5126	42	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 4.

All (42) 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:194:ARG:HD3	1:A:242:LEU:HD21	1.62	0.81
1:A:198:LEU:HD22	1:A:202:ILE:HD11	1.62	0.79
1:A:31:GLU:HG3	1:A:46:PHE:CE1	2.21	0.75
1:A:189[B]:GLN:NE2	5:A:501:HOH:O	2.17	0.72
1:B:79:PRO:HB3	1:B:348:LEU:HD22	1.71	0.70
1:A:352:GLN:HE22	4:A:403:GOL:H32	1.57	0.69
1:B:26:ASP:OD1	1:B:305[B]:ASP:OD1	2.11	0.69
1:A:30:ALA:HA	1:A:305[B]:ASP:OD1	1.92	0.68
1:A:189[A]:GLN:NE2	1:A:189[A]:GLN:HA	2.10	0.67
1:A:189[A]:GLN:HE21	1:A:189[A]:GLN:HA	1.68	0.57

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:MET:N	5:A:503:HOH:O	2.39	0.56
1:B:305[B]:ASP:HB3	5:B:535:HOH:O	2.06	0.54
1:A:19:LEU:HD11	1:A:345:LEU:HD12	1.91	0.53
1:A:31:GLU:HG3	1:A:46:PHE:CZ	2.44	0.53
1:A:305[B]:ASP:HB2	5:A:518:HOH:O	2.08	0.52
1:A:120:ASP:OD2	1:A:126:ARG:NH2	2.42	0.52
1:A:178:ILE:HG21	1:A:208:ILE:HD13	1.92	0.52
1:A:107:VAL:CG1	1:A:111:LYS:HE3	2.41	0.51
1:A:87:ALA:HB2	1:A:166:LEU:HD22	1.93	0.50
1:A:63:ASP:OD2	3:A:402:A3R:N4N	2.45	0.50
1:A:16:ARG:HE	1:A:321:MET:HG2	1.77	0.49
1:B:178:ILE:HG21	1:B:208:ILE:HD13	1.95	0.48
1:A:80:ASP:OD1	4:A:403:GOL:H31	2.12	0.48
1:B:289:ASN:HB3	1:B:319:TYR:CZ	2.49	0.48
1:B:24:LEU:HD23	1:B:332:CYS:SG	2.53	0.48
1:B:87:ALA:HB2	1:B:166:LEU:HD22	1.97	0.47
1:B:190:GLY:O	1:B:192:LEU:N	2.48	0.47
1:B:347:GLU:OE2	1:B:351:LYS:HE2	2.14	0.46
1:B:120:ASP:OD2	1:B:126:ARG:NH2	2.48	0.46
1:A:198:LEU:CD2	1:A:202:ILE:HD11	2.42	0.45
1:A:289:ASN:HB3	1:A:319:TYR:CZ	2.52	0.44
1:B:96:LYS:HE3	1:B:96:LYS:HB2	1.85	0.43
1:A:40:LEU:CD2	1:A:282:ASP:HB3	2.49	0.43
1:A:189[B]:GLN:HE21	1:A:192:LEU:CD2	2.33	0.42
1:B:266:VAL:HB	1:B:267:PRO:HD3	2.02	0.42
1:A:33:GLU:OE2	1:A:305[A]:ASP:OD1	2.38	0.41
1:B:306:THR:HA	1:B:309:THR:HG22	2.02	0.41
1:A:198:LEU:HD22	1:A:202:ILE:CD1	2.42	0.41
1:A:24:LEU:HD23	1:A:332:CYS:SG	2.61	0.41
1:A:306:THR:HA	1:A:309:THR:HG22	2.02	0.40
1:B:40:LEU:N	1:B:41:PRO:CD	2.84	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	329/350 (94%)	320 (97%)	8 (2%)	1 (0%)	41	30
1	B	331/350 (95%)	322 (97%)	8 (2%)	1 (0%)	41	30
All	All	660/700 (94%)	642 (97%)	16 (2%)	2 (0%)	41	30

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	191	GLU
1	A	191	GLU

### 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	271/283 (96%)	265 (98%)	6 (2%)	52	44
1	B	273/283 (96%)	268 (98%)	5 (2%)	59	53
All	All	544/566 (96%)	533 (98%)	11 (2%)	60	48

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	135	TYR
1	A	198	LEU
1	A	258	HIS
1	A	305[A]	ASP
1	A	305[B]	ASP
1	A	351	LYS
1	B	228	LYS
1	B	258	HIS
1	B	305[A]	ASP
1	B	305[B]	ASP
1	B	331	SER

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

Mol	Chain	Res	Type
1	A	154	GLN
1	A	185	HIS
1	A	246	ASN
1	A	352	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 for Mogul analysis.

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	$\# Z  > 2$
4	GOL	A	403	-	5,5,5	0.29	0	5,5,5	0.53	0
4	GOL	B	403	-	5,5,5	0.53	0	5,5,5	0.73	0
3	A3R	A	402	-	31,37,37	1.73	7 (22%)	33,56,56	2.38	6 (18%)
3	A3R	B	402	-	31,37,37	1.69	6 (19%)	33,56,56	1.59	6 (18%)

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
4	GOL	A	403	-	-	2/4/4/4	-
4	GOL	B	403	-	-	4/4/4/4	-
3	A3R	A	402	-	-	3/18/47/47	0/4/4/4
3	A3R	B	402	-	-	3/18/47/47	0/4/4/4

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	A3R	C2-N3	5.39	1.40	1.32
3	B	402	A3R	C2-N1	4.13	1.41	1.33
3	B	402	A3R	C2-N3	3.90	1.38	1.32
3	A	402	A3R	C2-N1	3.88	1.41	1.33
3	A	402	A3R	C3N-C2N	3.33	1.59	1.52
3	A	402	A3R	O2'-C2'	-3.19	1.35	1.43
3	B	402	A3R	C5N-C4N	3.14	1.58	1.51
3	B	402	A3R	C3'-C4'	-2.44	1.46	1.53
3	B	402	A3R	C6-C5	-2.38	1.34	1.43
3	B	402	A3R	C4-N3	-2.37	1.32	1.35
3	A	402	A3R	O3'-C3'	2.36	1.48	1.43
3	A	402	A3R	C6-C5	-2.19	1.35	1.43
3	A	402	A3R	PA-O2A	-2.02	1.45	1.55

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	402	A3R	C2N-C3N-C4N	-9.72	89.01	104.44
3	A	402	A3R	N3-C2-N1	-6.72	118.18	128.68
3	B	402	A3R	N3-C2-N1	-6.12	119.11	128.68
3	B	402	A3R	O2N-C2N-C3N	-2.75	103.22	110.03
3	A	402	A3R	C4-C5-N7	-2.41	106.88	109.40
3	A	402	A3R	O2N-C2N-C1N	-2.25	105.42	110.35
3	A	402	A3R	O2B-PB-O1B	2.24	123.33	112.24
3	B	402	A3R	C1'-N9-C4	-2.10	122.95	126.64
3	A	402	A3R	C3'-C2'-C1'	2.08	104.11	100.98
3	B	402	A3R	C2N-C3N-C4N	-2.08	101.14	104.44
3	B	402	A3R	O2A-PA-O1A	2.01	122.15	112.24
3	B	402	A3R	C3N-C4N-C5N	-2.00	109.73	113.26

There are no chirality outliers.

All (12) torsion outliers are listed below:

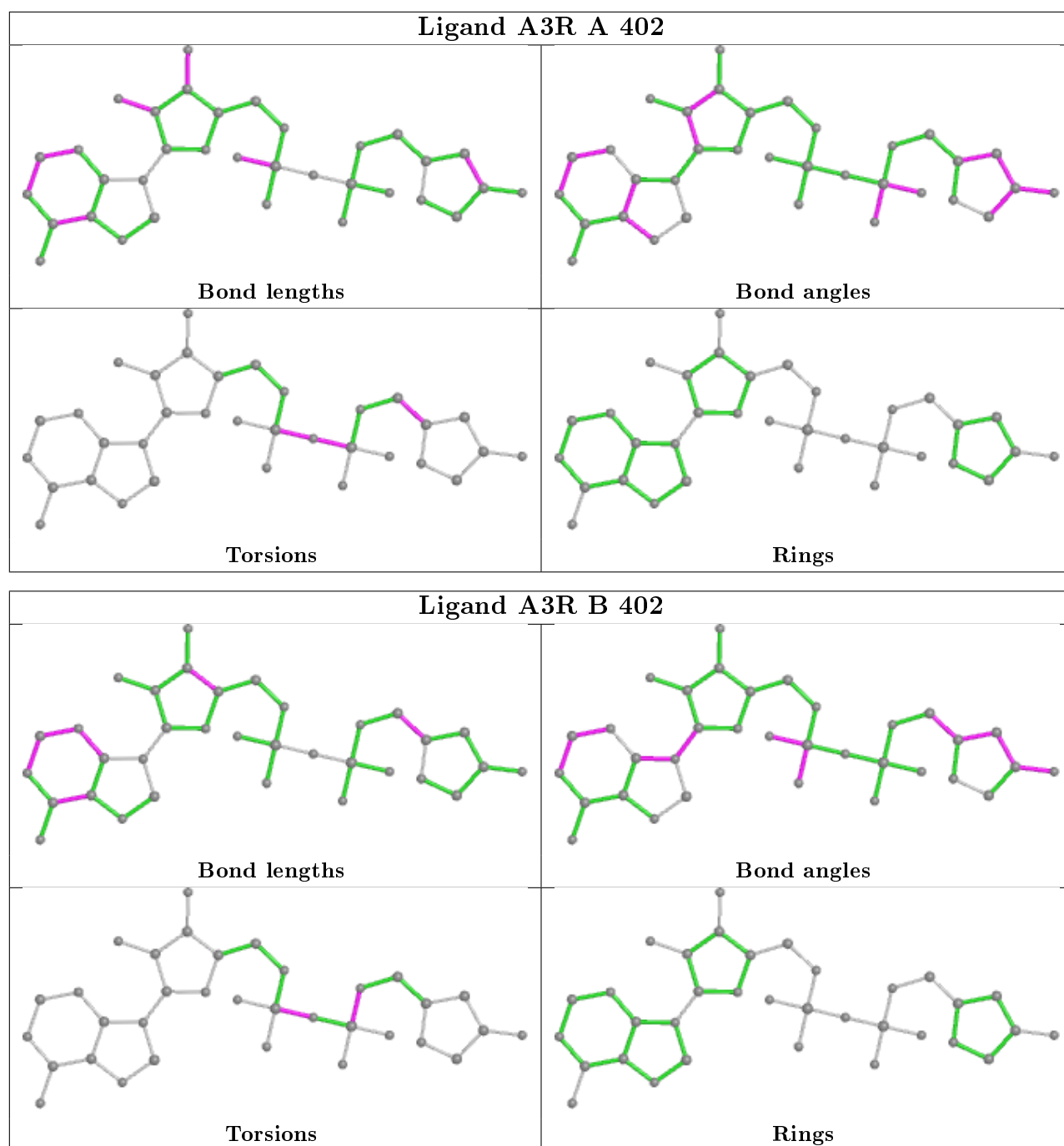
Mol	Chain	Res	Type	Atoms
3	A	402	A3R	PB-O3A-PA-O5'
3	B	402	A3R	C5N-O5N-PB-O1B
3	B	402	A3R	C5N-O5N-PB-O3A
3	B	402	A3R	PB-O3A-PA-O5'
4	A	403	GOL	C1-C2-C3-O3
4	B	403	GOL	O1-C1-C2-C3
4	A	403	GOL	O2-C2-C3-O3
4	B	403	GOL	C1-C2-C3-O3
4	B	403	GOL	O1-C1-C2-O2
4	B	403	GOL	O2-C2-C3-O3
3	A	402	A3R	PA-O3A-PB-O1B
3	A	402	A3R	N4N-C4N-C5N-O5N

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	403	GOL	2	0
3	A	402	A3R	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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	327/350 (93%)	0.83	51 (15%) 2 3	42, 65, 107, 122	0
1	B	328/350 (93%)	0.48	21 (6%) 19 28	35, 49, 78, 98	0
All	All	655/700 (93%)	0.65	72 (10%) 5 9	35, 55, 98, 122	0

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	232	CYS	6.0
1	A	233[A]	SER	5.2
1	A	37	ALA	5.2
1	A	116	PRO	5.1
1	B	229	LEU	5.0
1	A	229	LEU	5.0
1	A	230	PRO	4.7
1	A	114	LEU	4.6
1	A	235	LEU	4.6
1	A	250	ALA	4.5
1	A	245	SER	4.4
1	A	210	GLY	4.3
1	A	30	ALA	4.2
1	A	237	LYS	4.2
1	B	37	ALA	4.0
1	B	228	LYS	3.9
1	A	241	PHE	3.9
1	A	254	ASP	3.9
1	A	248	PRO	3.8
1	B	56	ALA	3.6
1	A	115	SER	3.6
1	A	238	ILE	3.5
1	A	240	GLU	3.5
1	A	252	ILE	3.5

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	B	116	PRO	3.5
1	B	54	LYS	3.4
1	A	202	ILE	3.4
1	A	32	PHE	3.3
1	A	236	LYS	3.3
1	B	36	ASP	3.3
1	B	233[A]	SER	3.3
1	A	255	GLU	3.3
1	A	52	LYS	3.1
1	A	304	THR	3.0
1	A	298	ILE	3.0
1	B	33[A]	GLU	2.9
1	B	236	LYS	2.9
1	A	244	SER	2.9
1	A	34	GLY	2.9
1	A	117	LYS	2.8
1	A	92[A]	GLU	2.8
1	B	52	LYS	2.7
1	A	36	ASP	2.7
1	A	247	VAL	2.7
1	A	251	ASP	2.6
1	B	232	CYS	2.6
1	A	209	GLU	2.6
1	B	245	SER	2.6
1	A	95	LYS	2.5
1	A	243	ALA	2.5
1	A	206	GLU	2.5
1	A	253	VAL	2.4
1	A	263	LEU	2.4
1	B	30	ALA	2.3
1	B	298	ILE	2.3
1	B	240	GLU	2.3
1	A	55	LYS	2.3
1	A	94	TYR	2.2
1	A	207	ARG	2.2
1	A	40	LEU	2.2
1	B	143	VAL	2.2
1	A	231	PHE	2.2
1	A	242	LEU	2.2
1	B	55	LYS	2.1
1	B	92[A]	GLU	2.1
1	A	111	LYS	2.1

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	A	38	VAL	2.1
1	A	246	ASN	2.1
1	B	210	GLY	2.1
1	A	191	GLU	2.0
1	B	255	GLU	2.0
1	A	239	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 carbohydrates in this entry.

## 6.4 Ligands [i](#)

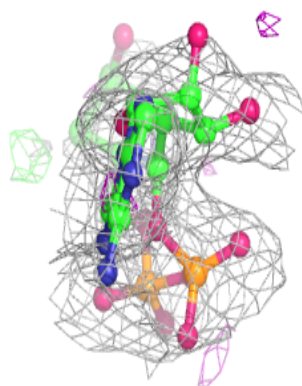
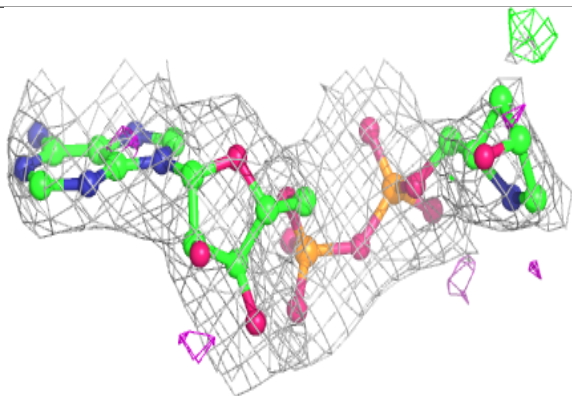
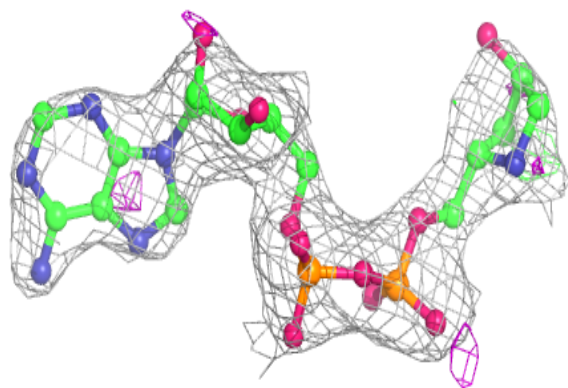
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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	GOL	B	403	6/6	0.81	0.26	89,90,94,96	0
4	GOL	A	403	6/6	0.86	0.24	81,84,85,87	0
2	MG	B	401	1/1	0.93	0.11	44,44,44,44	0
2	MG	A	401	1/1	0.96	0.08	62,62,62,62	0
3	A3R	A	402	34/34	0.96	0.12	61,72,84,87	0
3	A3R	B	402	34/34	0.96	0.14	45,48,55,56	0

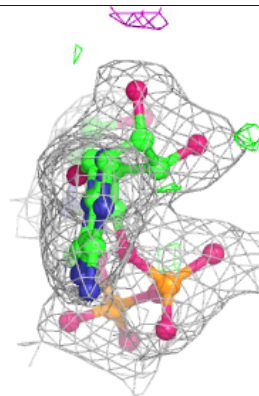
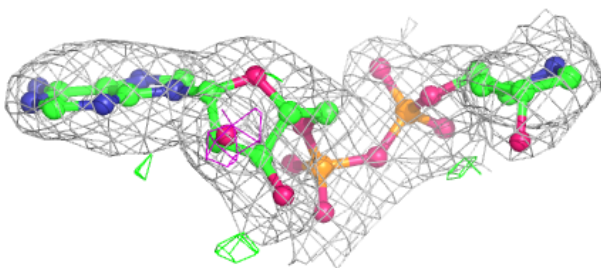
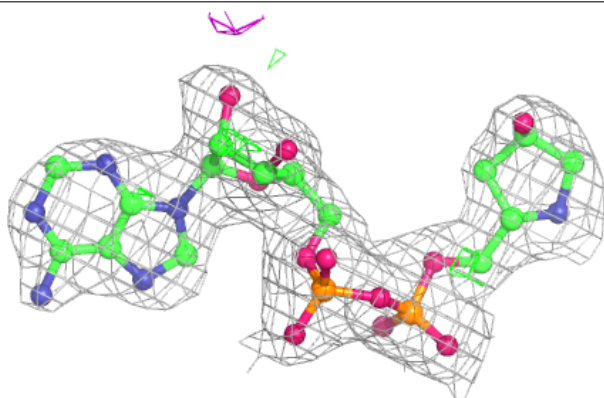
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

**Electron density around A3R A 402:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around A3R B 402:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



## 6.5 Other polymers

There are no such residues in this entry.