



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 03:38 pm GMT

PDB ID : 5DW8
Title : Crystal structure of 2'AMP bound SaIMPase-II
Authors : Dutta, A.; Bhattacharyya, S.; Das, A.K.
Deposited on : 2015-09-22
Resolution : 2.40 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

<http://wwpdb.org/validation/2016/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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

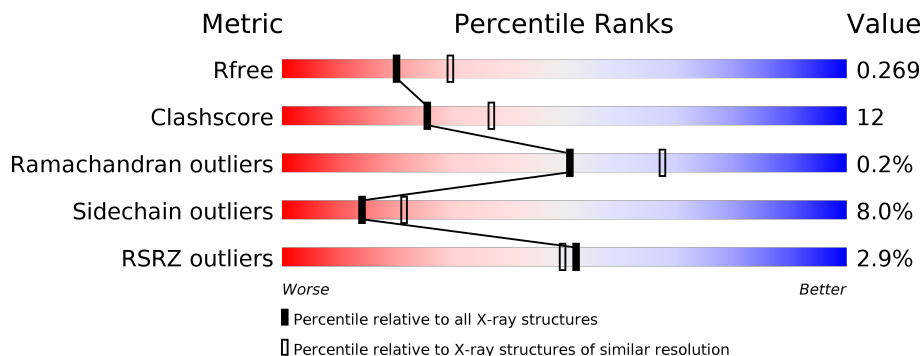
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	100719	3166 (2.40-2.40)
Clashscore	112137	3674 (2.40-2.40)
Ramachandran outliers	110173	3616 (2.40-2.40)
Sidechain outliers	110143	3617 (2.40-2.40)
RSRZ outliers	101464	3195 (2.40-2.40)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	283	<div> <div>3%</div> <div> <div></div> <div>70%</div> <div>20%</div> <div>8%</div> </div> </div>
1	B	283	<div> <div>2%</div> <div> <div></div> <div>73%</div> <div>14%</div> <div>8%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	2AM	A	302	-	-	-	X
3	2AM	B	303	-	-	-	X
4	GOL	A	304	-	-	X	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4192 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 Inositol monophosphatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	260	Total	C	N	O	S	0	1	0
			2014	1280	339	392	3			
1	B	259	Total	C	N	O	S	0	1	0
			2017	1285	340	388	4			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	HIS	-	expression tag	UNP A0A0C5I0W4
A	-6	HIS	-	expression tag	UNP A0A0C5I0W4
A	-5	HIS	-	expression tag	UNP A0A0C5I0W4
A	-4	HIS	-	expression tag	UNP A0A0C5I0W4
A	-3	HIS	-	expression tag	UNP A0A0C5I0W4
A	-2	HIS	-	expression tag	UNP A0A0C5I0W4
A	-1	GLY	-	expression tag	UNP A0A0C5I0W4
A	0	SER	-	expression tag	UNP A0A0C5I0W4
B	-7	HIS	-	expression tag	UNP A0A0C5I0W4
B	-6	HIS	-	expression tag	UNP A0A0C5I0W4
B	-5	HIS	-	expression tag	UNP A0A0C5I0W4
B	-4	HIS	-	expression tag	UNP A0A0C5I0W4
B	-3	HIS	-	expression tag	UNP A0A0C5I0W4
B	-2	HIS	-	expression tag	UNP A0A0C5I0W4
B	-1	GLY	-	expression tag	UNP A0A0C5I0W4
B	0	SER	-	expression tag	UNP A0A0C5I0W4

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Ca	0	0
			2	2		
2	A	2	Total	Ca	0	0
			2	2		

- # 2AM
-
- The chemical structure of 2-aminoadenosine (2AM) is shown. It consists of an adenine base (a purine ring system with an amino group at position 6) attached to a ribose sugar (a five-membered ring with hydroxyl groups at positions 2', 3', and 5'). The ribose sugar is linked to a phosphate group (a phosphorus atom double-bonded to one oxygen and single-bonded to three others, one of which is a hydroxyl group). The atoms are labeled with their respective symbols and positions: N1, N3, N6, C2, C4, C5, C6, C8, C9, N9, C1'(R), C2'(R), C3'(R), C4'(R), C5', O4', O5', O2', O3P, O2P, O3', O1P, and P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 23	C 10	N 5	O 7	P 1	0	0
3	B	1	Total 23	C 10	N 5	O 7	P 1	0	0

- # GOL
-
- The diagram shows the skeletal structure of 1,2,3-propanetriol (glycerol). The carbon backbone is represented by a zigzag line with three vertices labeled C1, C2, and C3 in green. C1 is on the left, C2 is in the middle, and C3 is on the right. Each carbon is bonded to a hydroxyl group (OH) in red. The OH group on C1 is labeled O1 in green below it. The OH group on C2 is labeled O2 in green below it. The OH group on C3 is labeled O3 in green to its right. The bonds between the carbons and the bonds to the hydroxyl groups are shown as thin grey lines.

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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	53	Total	O	0	0
			53	53		
5	B	52	Total	O	0	0
			52	52		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	49.01Å 105.69Å 53.31Å 90.00° 101.91° 90.00°	Depositor
Resolution (Å)	19.68 – 2.40 19.68 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.6 (19.68-2.40) 99.8 (19.68-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.94 (at 2.41Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.205 , 0.266 0.210 , 0.269	Depositor DCC
R_{free} test set	1064 reflections (5.41%)	DCC
Wilson B-factor (Å ²)	18.7	Xtriage
Anisotropy	0.230	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 41.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4192	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.81% 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: GOL, CA, 2AM

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	0.66	0/2061	0.82	1/2808 (0.0%)
1	B	0.71	0/2065	0.85	2/2808 (0.1%)
All	All	0.69	0/4126	0.83	3/5616 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	81	VAL	CB-CA-C	6.19	123.16	111.40
1	B	244	ARG	NE-CZ-NH1	5.54	123.07	120.30
1	A	216	LEU	CA-CB-CG	5.27	127.42	115.30

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	2014	0	1939	55	0
1	B	2017	0	1957	42	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	23	0	12	0	0
3	B	23	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	6	0	8	15	0
5	A	53	0	0	9	0
5	B	52	0	0	4	0
All	All	4192	0	3928	95	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 12.

All (95) 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:155:ASN:OD1	4:A:304:GOL:H31	1.79	0.82
1:A:155:ASN:OD1	4:A:304:GOL:C3	2.31	0.79
1:B:155:ASN:ND2	1:B:157:ASN:HD22	1.83	0.77
1:A:155:ASN:HD21	1:A:157:ASN:HD22	1.30	0.77
1:A:160:THR:HG22	1:B:169:LYS:HG3	1.66	0.77
1:B:248:GLN:NE2	5:B:401:HOH:O	2.21	0.74
1:A:40:ASN:O	1:A:41:VAL:HB	1.87	0.73
1:A:180:ALA:HB1	4:A:304:GOL:O3	1.88	0.72
1:B:217:TYR:HB3	5:B:412:HOH:O	1.88	0.72
1:A:155:ASN:HD22	1:A:157:ASN:H	1.38	0.71
1:A:155:ASN:ND2	1:A:157:ASN:HD22	1.91	0.69
1:A:176:ARG:HD3	5:A:439:HOH:O	1.94	0.68
1:A:65:GLU:OE1	1:A:66:GLU:OE1	2.12	0.67
1:A:155:ASN:HA	4:A:304:GOL:O3	1.93	0.67
1:A:180:ALA:HB1	4:A:304:GOL:C3	2.25	0.67
1:B:23:GLU:C	1:B:25:ASN:H	1.99	0.66
1:A:180:ALA:CB	4:A:304:GOL:O3	2.44	0.66
1:A:46:GLU:OE2	1:A:68:HIS:HD2	1.80	0.64
1:B:155:ASN:HD21	1:B:157:ASN:HD22	1.45	0.64
1:A:155:ASN:CA	4:A:304:GOL:O3	2.47	0.63
1:A:251:SER:HB2	5:A:402:HOH:O	1.99	0.61
1:A:203:ARG:HG2	1:A:269:ARG:NH1	2.15	0.61
1:A:182:GLY:HA2	4:A:304:GOL:O2	2.02	0.60
1:B:198:ALA:HB2	1:B:246:LEU:HD21	1.84	0.59
1:A:251:SER:CB	5:A:402:HOH:O	2.49	0.59
1:B:69:GLY:O	1:B:72:ILE:HG12	2.04	0.58
1:A:127:LYS:HE3	1:A:130:GLU:CD	2.25	0.57
1:A:40:ASN:O	1:A:41:VAL:CB	2.50	0.57
1:A:165:GLY:O	1:A:169:LYS:HD3	2.05	0.56
1:A:155:ASN:HB2	4:A:304:GOL:H31	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:251:SER:O	1:B:256:GLU:HB2	2.08	0.54
1:B:246:LEU:O	1:B:246:LEU:HG	2.06	0.54
1:A:180:ALA:HB1	4:A:304:GOL:H32	1.88	0.54
1:B:128:VAL:HB	1:B:217:TYR:OH	2.06	0.54
1:B:155:ASN:HD22	1:B:157:ASN:H	1.55	0.52
1:B:23:GLU:O	1:B:25:ASN:N	2.42	0.52
1:A:155:ASN:ND2	1:A:157:ASN:H	2.05	0.52
1:B:46[A]:GLU:HG3	1:B:50:PHE:CE2	2.45	0.52
1:B:46[A]:GLU:CG	1:B:50:PHE:CE2	2.94	0.51
1:B:152:ILE:HD12	1:B:171:ILE:HD11	1.93	0.51
1:A:81:VAL:CG2	1:A:207:TRP:HA	2.41	0.50
1:A:251:SER:O	1:A:256:GLU:HB2	2.11	0.50
1:B:170:GLU:HA	1:B:170:GLU:OE1	2.11	0.49
1:A:4:TYR:OH	1:A:125:HIS:ND1	2.41	0.49
1:A:57:TYR:HB3	1:A:60:HIS:CG	2.48	0.49
1:B:50:PHE:CD1	1:B:62:VAL:HG11	2.48	0.48
1:A:155:ASN:OD1	4:A:304:GOL:H32	2.14	0.48
1:A:169:LYS:HG3	1:B:160:THR:HG22	1.95	0.47
1:B:50:PHE:CE1	1:B:62:VAL:HG11	2.49	0.47
1:A:210:ALA:HB3	5:A:415:HOH:O	2.14	0.47
1:A:225:ASN:HB2	1:A:238:SER:OG	2.14	0.47
1:A:155:ASN:CB	4:A:304:GOL:H31	2.44	0.47
1:B:256:GLU:OE2	1:B:259:HIS:CD2	2.67	0.47
1:B:214:VAL:HA	1:B:217:TYR:CE1	2.50	0.47
1:A:74:THR:HB	1:A:233:ILE:HD12	1.97	0.46
1:A:151:ILE:HD12	1:B:93:HIS:CG	2.50	0.46
1:A:84:PRO:O	1:A:100:ILE:HG12	2.15	0.46
1:B:23:GLU:C	1:B:25:ASN:N	2.69	0.46
1:B:231:LEU:CD2	1:B:239:ILE:HD13	2.46	0.46
1:B:24:GLN:O	1:B:24:GLN:HG2	2.15	0.46
1:A:155:ASN:CG	4:A:304:GOL:H31	2.36	0.45
1:A:149:GLN:HG3	5:A:439:HOH:O	2.16	0.45
1:B:38:VAL:HG13	1:B:39:THR:N	2.32	0.45
1:A:127:LYS:HE3	1:A:130:GLU:OE1	2.18	0.44
1:B:125:HIS:CE1	1:B:133:TYR:HB2	2.52	0.44
1:B:152:ILE:CD1	1:B:171:ILE:HD11	2.48	0.44
1:B:205:GLN:NE2	5:B:405:HOH:O	2.50	0.44
1:A:168:PHE:O	1:A:172:VAL:HG23	2.17	0.44
1:B:191:VAL:HG21	1:B:241:VAL:HG23	1.99	0.44
1:B:59:ASN:HA	1:B:59:ASN:HD22	1.56	0.44
1:A:59:ASN:HB2	5:A:434:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:231:LEU:HD22	1:B:239:ILE:HD13	1.99	0.43
1:A:150:SER:HB3	1:A:197:GLU:HB2	2.00	0.43
1:A:81:VAL:HG22	1:A:207:TRP:HA	1.99	0.43
1:B:157:ASN:O	1:B:160:THR:OG1	2.23	0.43
1:B:72:ILE:HA	1:B:72:ILE:HD13	1.81	0.43
1:B:222:GLN:OE1	1:B:247:HIS:ND1	2.52	0.43
1:B:203:ARG:HB3	1:B:269:ARG:HD3	2.01	0.43
1:A:74:THR:HB	1:A:233:ILE:CD1	2.49	0.42
1:B:50:PHE:CE1	1:B:62:VAL:CG1	3.02	0.42
1:A:54:LEU:N	1:A:54:LEU:HD23	2.35	0.42
1:B:152:ILE:HG22	1:B:198:ALA:HB3	2.02	0.42
1:A:125:HIS:NE2	1:A:135:GLY:O	2.52	0.42
1:A:238:SER:O	1:A:239:ILE:HD13	2.19	0.42
1:A:158:TRP:HA	1:A:158:TRP:CE3	2.54	0.41
4:A:304:GOL:H11	5:A:431:HOH:O	2.20	0.41
1:A:82:VAL:HA	1:A:101:SER:O	2.20	0.41
1:A:155:ASN:HB2	4:A:304:GOL:C3	2.51	0.41
1:B:204:LEU:HG	1:B:208:ASP:HB2	2.03	0.41
1:A:155:ASN:HD22	1:A:157:ASN:N	2.14	0.41
1:B:60:HIS:CE1	1:B:78:THR:HG21	2.56	0.40
1:A:12:GLN:NE2	5:A:408:HOH:O	2.53	0.40
1:A:203:ARG:HB3	1:A:269:ARG:HD3	2.04	0.40
1:A:217:TYR:HE2	5:A:425:HOH:O	2.03	0.40
1:B:51:ASP:HB3	5:B:442:HOH:O	2.21	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	257/283 (91%)	245 (95%)	12 (5%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	256/283 (90%)	248 (97%)	7 (3%)	1 (0%)	38	54
All	All	513/566 (91%)	493 (96%)	19 (4%)	1 (0%)	51	67

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	24	GLN

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	213/239 (89%)	195 (92%)	18 (8%)	12	19
1	B	214/239 (90%)	198 (92%)	16 (8%)	16	24
All	All	427/478 (89%)	393 (92%)	34 (8%)	14	21

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	GLU
1	A	25	ASN
1	A	35	ASN
1	A	42	ASP
1	A	81	VAL
1	A	82	VAL
1	A	115	VAL
1	A	140	LYS
1	A	169	LYS
1	A	170	GLU
1	A	171	ILE
1	A	174	ASP
1	A	203	ARG
1	A	213	LEU
1	A	222	GLN

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Mol	Chain	Res	Type
1	A	244	ARG
1	A	246	LEU
1	A	251	SER
1	B	25	ASN
1	B	59	ASN
1	B	72	ILE
1	B	81	VAL
1	B	115	VAL
1	B	134	ARG
1	B	146	ASN
1	B	170	GLU
1	B	171	ILE
1	B	174	ASP
1	B	203	ARG
1	B	213	LEU
1	B	244	ARG
1	B	246	LEU
1	B	251	SER
1	B	264	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	GLN
1	A	20	GLN
1	A	24	GLN
1	A	68	HIS
1	A	70	HIS
1	A	93	HIS
1	A	157	ASN
1	A	222	GLN
1	A	259	HIS
1	A	268	GLN
1	B	59	ASN
1	B	93	HIS
1	B	146	ASN
1	B	155	ASN
1	B	205	GLN
1	B	259	HIS
1	B	268	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 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$
3	2AM	A	302	2	21,25,25	2.38	7 (33%)	25,38,38	2.76	4 (16%)
4	GOL	A	304	-	5,5,5	1.23	1 (20%)	5,5,5	1.52	1 (20%)
3	2AM	B	303	2	21,25,25	2.34	6 (28%)	25,38,38	2.71	3 (12%)

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	2AM	A	302	2	-	0/7/27/27	0/3/3/3
4	GOL	A	304	-	-	0/4/4/4	0/0/0/0
3	2AM	B	303	2	-	0/7/27/27	0/3/3/3

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	302	2AM	C5-C4	-2.42	1.35	1.40
3	B	303	2AM	C5-C4	-2.29	1.35	1.40
4	A	304	GOL	C3-C2	-2.03	1.44	1.52
3	A	302	2AM	P-O2P	2.05	1.63	1.54
3	B	303	2AM	O4'-C1'	2.42	1.44	1.41
3	A	302	2AM	P-O1P	3.08	1.61	1.50
3	B	303	2AM	C2-N1	3.46	1.40	1.33
3	B	303	2AM	P-O1P	3.68	1.63	1.50
3	A	302	2AM	C2-N1	3.69	1.40	1.33
3	B	303	2AM	P-O2'	4.40	1.67	1.59
3	A	302	2AM	P-O2'	4.69	1.67	1.59
3	A	302	2AM	O4'-C1'	4.70	1.47	1.41
3	A	302	2AM	C2-N3	5.65	1.41	1.32
3	B	303	2AM	C2-N3	6.46	1.42	1.32

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	302	2AM	N3-C2-N1	-12.26	118.18	128.86
3	B	303	2AM	N3-C2-N1	-11.92	118.48	128.86
4	A	304	GOL	O2-C2-C3	-3.22	93.61	108.84
3	B	303	2AM	C5-C6-N6	-2.20	115.98	120.47
3	A	302	2AM	C5-C6-N6	-2.16	116.06	120.47
3	A	302	2AM	O4'-C1'-C2'	-2.10	102.92	106.59
3	A	302	2AM	O3P-P-O2'	2.18	115.90	106.00
3	B	303	2AM	O4'-C4'-C5'	2.26	113.99	109.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	304	GOL	15	0

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	260/283 (91%)	0.17	9 (3%)	44 43	10, 18, 40, 56	2 (0%)
1	B	259/283 (91%)	0.04	6 (2%)	61 58	9, 16, 30, 41	2 (0%)
All	All	519/566 (91%)	0.10	15 (2%)	52 50	9, 17, 35, 56	4 (0%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	5.3
1	A	267	GLU	3.9
1	A	24	GLN	3.4
1	A	23	GLU	2.8
1	A	36	ASP	2.8
1	A	25	ASN	2.7
1	A	35	ASN	2.7
1	B	39	THR	2.5
1	B	25	ASN	2.4
1	B	24	GLN	2.4
1	B	270	PHE	2.3
1	B	67	GLY	2.2
1	A	40	ASN	2.1
1	B	217	TYR	2.1
1	A	268	GLN	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
4	GOL	A	304	6/6	0.86	0.32	8.62	21,22,24,26	0
3	2AM	B	303	23/23	0.82	0.26	4.82	28,40,61,63	0
3	2AM	A	302	23/23	0.84	0.28	3.49	37,45,64,65	0
2	CA	A	303	1/1	0.90	0.17	0.85	70,70,70,70	0
2	CA	B	301	1/1	0.99	0.03	-3.85	22,22,22,22	0
2	CA	A	301	1/1	0.93	0.07	-4.30	45,45,45,45	0
2	CA	B	302	1/1	0.97	0.05	-4.38	31,31,31,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.