

# Full wwPDB X-ray Structure Validation Report (i)

#### Oct 9, 2023 – 10:56 PM EDT

PDB ID : 7MQT

Title: N-Acetylmannosamine-6-phosphate 2-epimerase from Staphylococcus aureus

(strain MRSA USA300) with 5-deoxy substrate bound

Authors : Currie, M.J.; Dobson, R.C.J.

Deposited on : 2021-05-06

Resolution : 1.88 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 (i) 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 (1)) 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 : 2.35.1

buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$ 

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

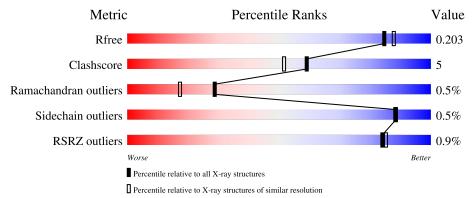
Validation Pipeline (wwPDB-VP) : 2.35.1

## 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X- $RAY\ DIFFRACTION$ 

The reported resolution of this entry is 1.88 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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries},{\rm resolution\ range}(\mathring{\rm A})) \end{array}$
$R_{free}$	130704	9470 (1.90-1.86)
Clashscore	141614	10282 (1.90-1.86)
Ramachandran outliers	138981	10152 (1.90-1.86)
Sidechain outliers	138945	10152 (1.90-1.86)
RSRZ outliers	127900	9303 (1.90-1.86)

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 <=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	AAA	222	91%	9%				
1	BBB	222	91%	9%				



# 2 Entry composition (i)

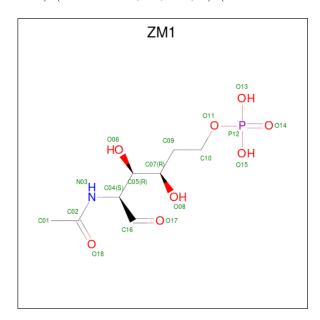
There are 5 unique types of molecules in this entry. The entry contains 4279 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 N-acetylmannosamine-6-phosphate 2-epimerase.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	AAA	222	Total 1808	C 1148	N 297	O 353	S 10	0	13	0
1	BBB	222	Total 1826	C 1156	N 299	O 360	S 11	0	16	0

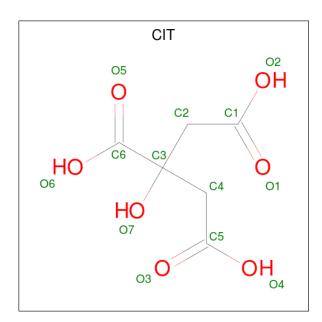
• Molecule 2 is 2-acetamido-2,5-dideoxy-6-O-phosphono-D-lyxo-hexose (three-letter code: ZM1) (formula: C<sub>8</sub>H<sub>16</sub>NO<sub>8</sub>P) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf			
2	AAA	1	Total	С	N	О	Р	0	1	
	AAA	1	18	8	1	8	1	0	1	
2	BBB	1	Total	С	N	О	Р	0	1	
2	DDD	1	18	8	1	8	1		1	

• Molecule 3 is CITRIC ACID (three-letter code: CIT) (formula: C<sub>6</sub>H<sub>8</sub>O<sub>7</sub>).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	AAA	1	Total C O 13 6 7	0	1
3	BBB	1	Total C O 13 6 7	0	1

• Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

I	Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
	4	AAA	1	Total Cl 1 1	0	0
	4	BBB	1	Total Cl 1 1	0	0

• Molecule 5 is water.

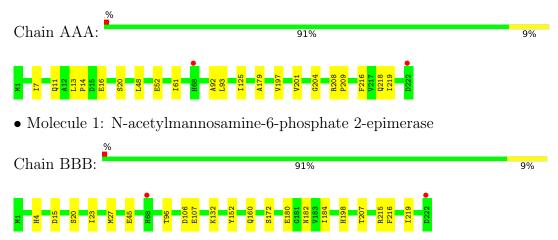
$\mathbf{Mol}$	Chain	Residues	${f Atoms}$	ZeroOcc	AltConf
5	AAA	288	Total O 288 288	0	0
5	BBB	293	Total O 293 293	0	0



# 3 Residue-property plots (i)

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: N-acetylmannosamine-6-phosphate 2-epimerase





# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants	76.44Å 173.45Å 44.04Å	Donositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.11 - 1.88	Depositor
Resolution (A)	46.11 - 1.88	EDS
% Data completeness	99.7 (46.11-1.88)	Depositor
(in resolution range)	99.7 (46.11-1.88)	EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.39 (at 1.88Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
D D.	0.154 , 0.197	Depositor
$R, R_{free}$	0.164 , 0.203	DCC
$R_{free}$ test set	2470 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	20.6	Xtriage
Anisotropy	0.174	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.37, 44.5	EDS
L-test for twinning <sup>2</sup>	$ < L >=0.50, < L^2>=0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	4279	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.96% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>Theoretical values of <|L|>,  $<L^2>$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



<sup>&</sup>lt;sup>1</sup>Intensities estimated from amplitudes.

# 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: CL, CIT, ZM1

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		Bo	nd lengths	Bond angles		
IVIOI	Chain	RMSZ	# Z  > 5	RMSZ	# Z >5	
1	AAA	0.75	1/1857 (0.1%)	0.87	0/2513	
1	BBB	0.72	0/1877	0.83	0/2542	
All	All	0.73	1/3734 (0.0%)	0.85	0/5055	

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\text{\AA})$	$\operatorname{Ideal}( ext{\AA})$
1	AAA	16	GLU	CD-OE1	5.02	1.31	1.25

There are no bond angle outliers.

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	AAA	1808	0	1836	12	0
1	BBB	1826	0	1845	19	0
2	AAA	18	0	0	0	0
2	BBB	18	0	0	4	0
3	AAA	13	0	5	3	0
3	BBB	13	0	5	0	0
4	AAA	1	0	0	0	0
4	BBB	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	$\mathbf{H}(\mathbf{added})$	Clashes	Symm-Clashes
5	AAA	288	0	0	1	2
5	BBB	293	0	0	11	2
All	All	4279	0	3691	35	2

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 5.

All (35) 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:BBB:15[A]:ASP:OD1	5:BBB:405:HOH:O	1.84	0.94
3:AAA:302[A]:CIT:C6	3:AAA:302[A]:CIT:O3	2.28	0.81
1:BBB:215[B]:ARG:HD3	5:BBB:464:HOH:O	1.89	0.71
2:BBB:301[B]:ZM1:O15	5:BBB:407:HOH:O	2.08	0.71
1:BBB:172[A]:SER:OG	5:BBB:408:HOH:O	2.11	0.69
1:AAA:48:LEU:O	1:AAA:52[B]:GLU:HG3	1.96	0.66
1:BBB:45[B]:GLU:OE1	5:BBB:401:HOH:O	0.58	0.58
1:BBB:160:GLN:OE1	5:BBB:410:HOH:O	2.17	0.57
1:BBB:215[B]:ARG:CD	5:BBB:464:HOH:O	2.51	0.55
2:BBB:301[B]:ZM1:P12	5:BBB:407:HOH:O	2.67	0.53
1:AAA:11:GLN:HE22	3:AAA:302[A]:CIT:C1	2.22	0.53
1:AAA:209:PRO:HG3	1:BBB:27[A]:MET:SD	2.49	0.53
1:BBB:180:GLU:OE2	2:BBB:301[B]:ZM1:O08	2.26	0.52
1:AAA:216:PHE:O	1:AAA:219[A]:ILE:HG22	2.11	0.51
1:AAA:7[B]:ILE:O	1:AAA:201:VAL:HA	2.14	0.47
1:BBB:132:LYS:HE2	1:BBB:172[A]:SER:HB3	1.98	0.46
1:BBB:27[B]:MET:CE	1:BBB:207:THR:HA	2.45	0.46
1:BBB:107[A]:GLU:HG2	5:BBB:454:HOH:O	2.17	0.45
1:BBB:216:PHE:O	1:BBB:219[B]:ILE:HG22	2.18	0.43
1:BBB:23:ILE:O	1:BBB:27[B]:MET:HG2	2.18	0.43
1:AAA:204:GLY:HA2	1:AAA:208:ARG:HB2	2.01	0.43
1:AAA:13:LEU:HB3	1:AAA:14:PRO:HD2	2.01	0.43
1:BBB:215[C]:ARG:HD3	5:BBB:409:HOH:O	2.18	0.43
3:AAA:302[A]:CIT:O3	3:AAA:302[A]:CIT:O5	2.37	0.42
1:BBB:182[B]:ASN:OD1	1:BBB:184[B]:ILE:HD12	2.19	0.42
1:AAA:61:ILE:CG2	1:AAA:92:ALA:HB3	2.49	0.42
1:AAA:218:GLN:NE2	5:AAA:425:HOH:O	2.53	0.42
1:AAA:7[A]:ILE:HB	1:AAA:201:VAL:HG22	2.02	0.41
1:BBB:45[A]:GLU:CD	5:BBB:498:HOH:O	2.58	0.41
2:BBB:301[B]:ZM1:O06	2:BBB:301[B]:ZM1:O11	2.37	0.41
1:BBB:4:HIS:HA	1:BBB:198:HIS:O	2.21	0.41

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Atom-1	Atom-2	$egin{aligned} &  ext{Interatomic} \ &  ext{distance} \ &  ext{(Å)} \end{aligned}$	Clash overlap (Å)
1:AAA:179:ALA:HB2	1:AAA:197:VAL:HG11	2.03	0.40
1:BBB:96:THR:HB	1:BBB:152:TYR:OH	2.21	0.40
1:BBB:132:LYS:HE2	1:BBB:172[B]:SER:HB2	2.03	0.40
1:AAA:125:ILE:HD12	1:AAA:125:ILE:C	2.42	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
5:AAA:651:HOH:O	5:BBB:665:HOH:O[1_556]	2.06	0.14
5:AAA:615:HOH:O	5:BBB:533:HOH:O[4_593]	2.14	0.06

### 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	Perce	$\mathbf{ntiles}$
1	AAA	$233/222\ (105\%)$	226 (97%)	6 (3%)	1 (0%)	34	22
1	BBB	$236/222\ (106\%)$	229 (97%)	6 (2%)	1 (0%)	34	22
All	All	469/444~(106%)	455 (97%)	12 (3%)	2 (0%)	29	22

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	AAA	20	SER
1	BBB	20	SER

#### 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	P	erce	ntiles
1	AAA	203/190 (107%)	202 (100%)	1 (0%)		88	88
1	BBB	206/190~(108%)	205 (100%)	1 (0%)		88	88
All	All	409/380 (108%)	407 (100%)	2 (0%)		88	88

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	93	LEU
1	BBB	106	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

#### 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)

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	Trunc	Chain	Res Link		Во	ond leng	ths	В	ond ang	les
IVIOI	Type	Chain	nes	Link	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	CIT	AAA	302[A]	-	12,12,12	1.19	1 (8%)	17,17,17	1.42	2 (11%)
2	ZM1	BBB	301[B]	-	16,17,17	1.21	1 (6%)	17,23,23	1.22	3 (17%)
3	CIT	BBB	302[A]	-	12,12,12	1.23	1 (8%)	17,17,17	1.47	4 (23%)
2	ZM1	AAA	301[B]	-	16,17,17	1.05	1 (6%)	17,23,23	1.30	1 (5%)

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	CIT	AAA	302[A]	-	-	9/16/16/16	-
2	ZM1	BBB	301[B]	-	-	8/19/21/21	-
3	CIT	BBB	302[A]	-	-	8/16/16/16	-
2	ZM1	AAA	301[B]	-	-	8/19/21/21	-

#### All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(A)	Ideal(A)
2	BBB	301[B]	ZM1	C02-N03	3.61	1.46	1.34
3	BBB	302[A]	CIT	C3-C6	3.02	1.56	1.53
2	AAA	301[B]	ZM1	C02-N03	2.61	1.43	1.34
3	AAA	302[A]	CIT	C3-C6	2.07	1.55	1.53

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}(^{o})$
2	AAA	301[B]	ZM1	C04-N03-C02	-4.09	118.65	123.12
3	AAA	302[A]	CIT	O5-C6-C3	-3.76	116.93	122.25
3	BBB	302[A]	CIT	O5-C6-C3	-2.88	118.17	122.25
3	BBB	302[A]	CIT	O7-C3-C6	2.60	112.51	108.86
3	AAA	302[A]	CIT	O6-C6-C3	2.57	117.52	113.05
3	BBB	302[A]	CIT	O6-C6-C3	2.46	117.32	113.05
2	BBB	301[B]	ZM1	O18-C02-C01	-2.25	117.88	122.06
2	BBB	301[B]	ZM1	C04-N03-C02	-2.24	120.67	123.12
3	BBB	302[A]	CIT	O3-C5-C4	-2.20	116.52	122.94
2	BBB	301[B]	ZM1	P12-O11-C10	-2.02	112.73	118.30

There are no chirality outliers.

All (33) torsion outliers are listed below:



Mol	Chain	Res	Type	Atoms
2	AAA	301[B]	ZM1	N03-C04-C05-O06
2	AAA	301[B]	ZM1	C16-C04-C05-C07
2	AAA	301[B]	ZM1	O08-C07-C09-C10
2	AAA	301[B]	ZM1	C10-O11-P12-O14
2	BBB	301[B]	ZM1	N03-C04-C05-O06
2	BBB	301[B]	ZM1	N03-C04-C05-C07
2	BBB	301[B]	ZM1	C16-C04-C05-O06
2	BBB	301[B]	ZM1	C16-C04-C05-C07
2	BBB	301[B]	ZM1	O08-C07-C09-C10
2	BBB	301[B]	ZM1	C10-O11-P12-O13
2	BBB	301[B]	ZM1	C10-O11-P12-O15
3	AAA	302[A]	CIT	C2-C3-C6-O5
3	AAA	302[A]	CIT	C2-C3-C6-O6
3	AAA	302[A]	CIT	O7-C3-C6-O5
3	AAA	302[A]	CIT	O7-C3-C6-O6
3	AAA	302[A]	CIT	O7-C3-C4-C5
3	AAA	302[A]	CIT	C6-C3-C4-C5
2	AAA	301[B]	ZM1	C16-C04-C05-O06
3	BBB	302[A]	CIT	C2-C3-C6-O5
3	AAA	302[A]	CIT	C2-C3-C4-C5
2	AAA	301[B]	ZM1	C05-C07-C09-C10
3	BBB	302[A]	CIT	O7-C3-C6-O5
3	BBB	302[A]	CIT	O7-C3-C6-O6
3	AAA	302[A]	CIT	C4-C3-C6-O6
3	BBB	302[A]	CIT	C2-C3-C6-O6
2	BBB	301[B]	ZM1	C05-C07-C09-C10
2	AAA	301[B]	ZM1	N03-C04-C05-C07
3	AAA	302[A]	CIT	C4-C3-C6-O5
3	BBB	302[A]	CIT	C4-C3-C6-O6
3	BBB	302[A]	CIT	C4-C3-C6-O5
3	BBB	302[A]	CIT	O2-C1-C2-C3
3	BBB	302[A]	CIT	O1-C1-C2-C3
2	AAA	301[B]	ZM1	C10-O11-P12-O15

There are no ring outliers.

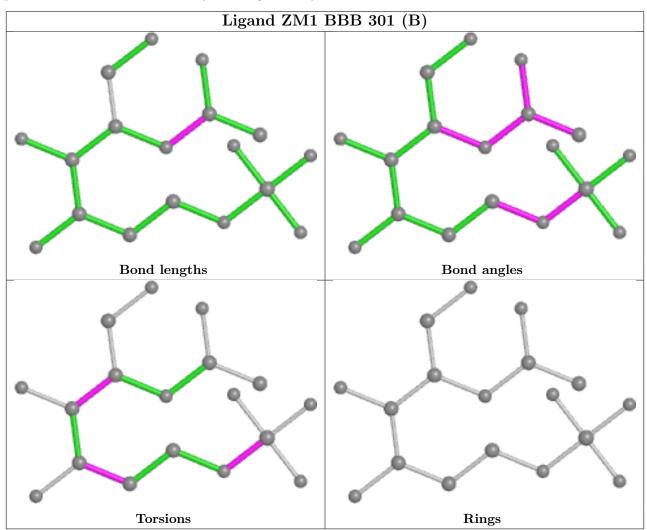
2 monomers are involved in 7 short contacts:

	Mol	Chain	Res	Type	Clashes	Symm-Clashes
	3		302[A]		3	0
Ī	2	BBB	301[B]	ZM1	4	0

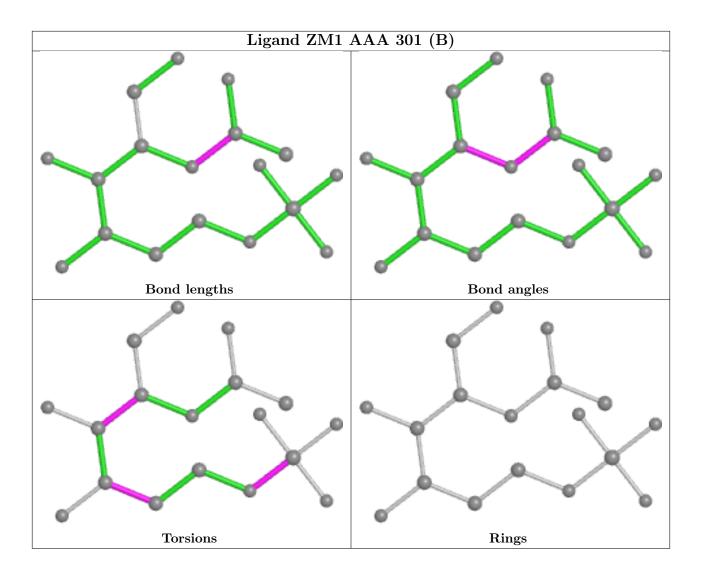
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In



addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 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 (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,  $95^{th}$  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	$\begin{array}{c c} \textbf{Analysed} &  <\!\!\text{RSRZ}\!\!>\!  & \#\!\!\text{RSRZ}\!\!>\!\!2 \end{array}$		2	$\mathbf{OWAB}(\mathbf{\mathring{A}}^2)$	Q < 0.9	
1	AAA	222/222 (100%)	-0.32	2 (0%) 84 8	85	15, 20, 32, 56	0
1	BBB	$222/222 \ (100\%)$	-0.32	2 (0%) 84 8	85	14, 19, 31, 69	0
All	All	444/444 (100%)	-0.32	4 (0%) 84 8	85	14, 20, 32, 69	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	BBB	68	HIS	3.5
1	AAA	68	HIS	2.5
1	AAA	222	ASP	2.4
1	BBB	222	ASP	2.2

## 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,  $95^{th}$  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	$\operatorname{Res}$	Atoms	RSCC	RSR	$  $ B-factors $(A^2)$	Q<0.9
3	CIT	AAA	302[A]	13/13	0.92	0.14	26,29,32,32	13

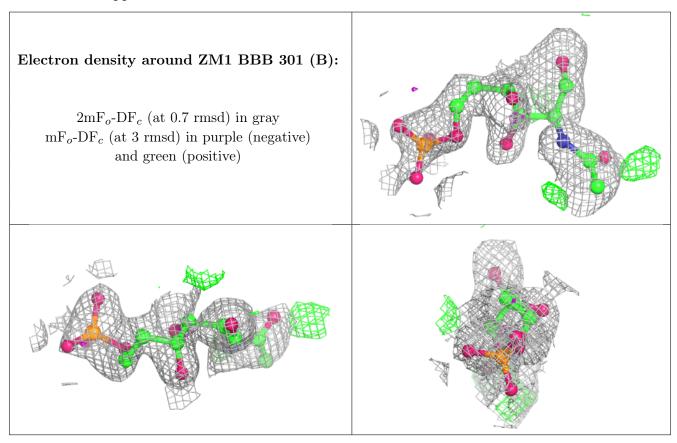
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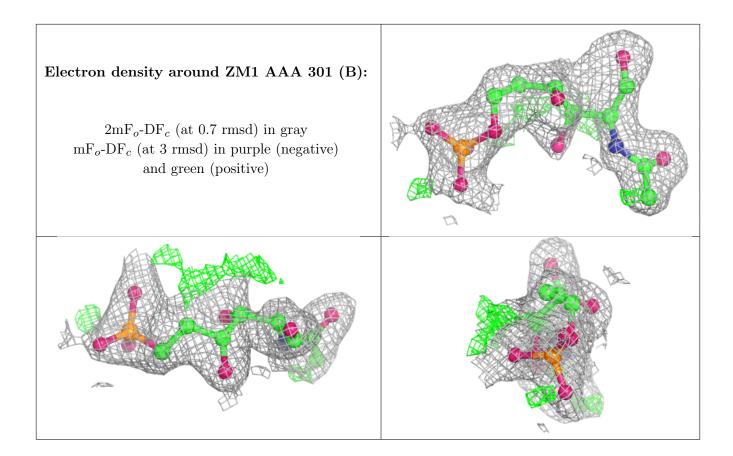
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	${f B-factors}({f \AA}^2)$	Q < 0.9
2	ZM1	BBB	301[B]	18/18	0.93	0.15	30,36,41,42	18
2	ZM1	AAA	301[B]	18/18	0.93	0.13	21,33,41,48	18
3	CIT	BBB	302[A]	13/13	0.94	0.12	17,19,22,24	13
4	CL	AAA	303	1/1	1.00	0.08	20,20,20,20	0
4	CL	BBB	303	1/1	1.00	0.08	23,23,23,23	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.







# 6.5 Other polymers (i)

There are no such residues in this entry.

