

Full wwPDB X-ray Structure Validation Report (i)

Aug 10, 2023 – 06:06 PM EDT

43 of Campylobacter
4

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:

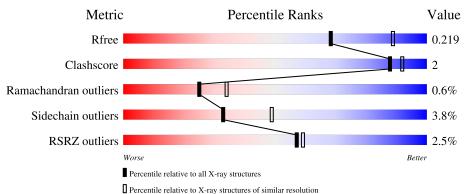
MolProbity	:	4.02b-467 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)		
EDS	:	2.35
buster-report	:	1.1.7(2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.38 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	5509(2.40-2.36)
Clashscore	141614	6082(2.40-2.36)
Ramachandran outliers	138981	5973 (2.40-2.36)
Sidechain outliers	138945	5975 (2.40-2.36)
RSRZ outliers	127900	5397 (2.40-2.36)

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	А	544	2% 8 6%	7% • 6%



$8 \mathrm{GFM}$

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 4463 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 Lytic transglycosylase domain-containing protein.

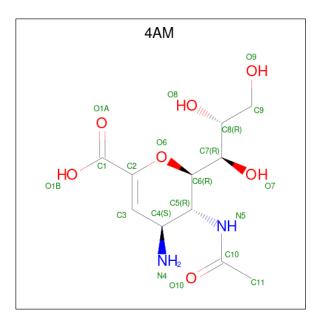
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	А	512	Total 4247	C 2755	N 697	O 780	S 15	0	1	0

Chain	Residue	Modelled	Actual	Comment	Reference
А	-2	MET	-	expression tag	UNP A0A3I4HTM2
A	-1	GLY	-	expression tag	UNP A0A3I4HTM2
А	0	SER	-	expression tag	UNP A0A3I4HTM2
А	1	SER	-	expression tag	UNP A0A3I4HTM2
A	2	HIS	-	expression tag	UNP A0A3I4HTM2
А	3	HIS	-	expression tag	UNP A0A3I4HTM2
А	4	HIS	-	expression tag	UNP A0A3I4HTM2
A	5	HIS	-	expression tag	UNP A0A3I4HTM2
А	6	HIS	-	expression tag	UNP A0A3I4HTM2
A	7	HIS	-	expression tag	UNP A0A3I4HTM2
А	8	SER	-	expression tag	UNP A0A3I4HTM2
A	9	SER	-	expression tag	UNP A0A3I4HTM2
А	10	GLY	-	expression tag	UNP A0A3I4HTM2
А	11	LEU	-	expression tag	UNP A0A3I4HTM2
А	12	VAL	-	expression tag	UNP A0A3I4HTM2
А	13	PRO	-	expression tag	UNP A0A3I4HTM2
А	14	ARG	-	expression tag	UNP A0A3I4HTM2
А	15	GLY	-	expression tag	UNP A0A3I4HTM2
А	16	SER	-	expression tag	UNP A0A3I4HTM2
А	17	HIS	-	expression tag	UNP A0A3I4HTM2
А	18	MET	-	expression tag	UNP A0A3I4HTM2

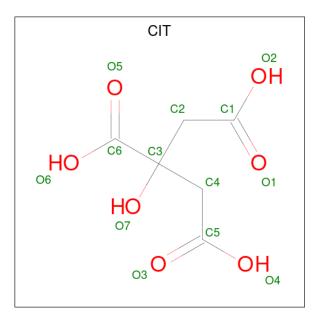
There are 21 discrepancies between the modelled and reference sequences:

• Molecule 2 is 4-AMINO-2-DEOXY-2,3-DEHYDRO-N-ACETYL-NEURAMINIC ACID (three-letter code: 4AM) (formula: C₁₁H₁₈N₂O₇) (labeled as "Ligand of Interest" by depositor).





Mol	Chain	Residues	A	Atoms			ZeroOcc	AltConf
2	А	1	Total 20	C 11	N 2	O 7	0	0



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

• Molecule 4 is water.



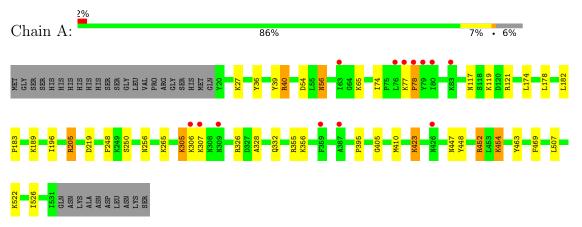
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	183	Total O 183 183	0	0



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: Lytic transglycosylase domain-containing protein





4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 2 3	Depositor
Cell constants	178.06Å 178.06 Å 178.06 Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	-
Resolution (Å)	29.68 - 2.38	Depositor
Resolution (A)	29.68 - 2.38	EDS
% Data completeness	99.8 (29.68-2.38)	Depositor
(in resolution range)	99.9(29.68-2.38)	EDS
R _{merge}	0.09	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.11 (at 2.36Å)	Xtriage
Refinement program	REFMAC 5.8.0349	Depositor
D D	0.174 , 0.214	Depositor
R, R_{free}	0.182 , 0.219	DCC
R_{free} test set	1932 reflections (5.11%)	wwPDB-VP
Wilson B-factor $(Å^2)$	44.7	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.32 , 30.2	EDS
L-test for twinning ²	$< L >=0.51, < L^2>=0.34$	Xtriage
Estimated twinning fraction	0.016 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4463	wwPDB-VP
Average B, all atoms $(Å^2)$	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 1.94% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities 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: CIT, $4\mathrm{AM}$

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	Bo	nd angles
	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5
1	А	0.45	0/4351	0.80	3/5871~(0.1%)

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	А	0	4

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	121	ARG	NE-CZ-NH2	-7.20	116.70	120.30
1	А	40	ARG	NE-CZ-NH2	-7.14	116.73	120.30
1	А	40	ARG	NE-CZ-NH1	5.54	123.07	120.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	А	326	ARG	Sidechain
1	А	355	ARG	Sidechain
1	А	452	ARG	Sidechain
1	А	54	ASP	Peptide



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	А	4247	0	4258	17	0
2	А	20	0	17	1	0
3	А	13	0	5	0	0
4	А	183	0	0	1	0
All	All	4463	0	4280	17	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 2.

All (17) 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:410[B]:MET:HE2	4:A:871:HOH:O	1.73	0.88
1:A:395:PRO:O	1:A:405:GLY:HA2	2.00	0.61
1:A:174:LEU:HD22	1:A:178:LEU:HD23	1.86	0.57
1:A:507:LEU:HD11	1:A:526:ILE:HG21	1.85	0.57
1:A:448:TYR:CE1	1:A:452:ARG:CZ	2.96	0.49
1:A:39:TYR:CE1	1:A:65:LYS:HG3	2.48	0.49
1:A:189:LYS:HD2	1:A:219:ASP:HB2	1.96	0.47
1:A:74:ILE:O	1:A:74:ILE:HG22	2.15	0.45
1:A:36:TYR:CZ	1:A:40:ARG:HD2	2.51	0.45
1:A:463:TYR:CE1	2:A:601:4AM:H112	2.52	0.44
1:A:77:LYS:N	1:A:78:PRO:CD	2.81	0.44
1:A:328:ALA:HB1	1:A:332:GLN:HB2	2.00	0.43
1:A:77:LYS:O	1:A:78:PRO:O	2.36	0.43
1:A:56:ASN:HD22	1:A:56:ASN:H	1.68	0.41
1:A:196:ILE:HA	1:A:205:ARG:HD2	2.03	0.41
1:A:182:LEU:N	1:A:183:PRO:CD	2.84	0.41
1:A:305:LYS:HD3	1:A:305:LYS:HA	1.75	0.41

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	А	511/544~(94%)	499 (98%)	9(2%)	3~(1%)	25 34

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	423	LYS
1	А	78	PRO
1	А	454	LYS

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	А	469/496~(95%)	451 (96%)	18 (4%)	33 49	

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	27	LYS
1	А	56	ASN
1	А	117	ASN
1	А	119	LYS
1	А	205	ARG
1	А	248	PHE
1	А	250	SER
1	А	256	ASN

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		-	Jus puye
Mol	Chain	\mathbf{Res}	Type
1	А	265	LYS
1	А	305	LYS
1	А	306	LYS
1	А	307	LYS
1	А	356	LYS
1	А	423	LYS
1	А	447	ASN
1	А	454	LYS
1	А	469	PHE
1	А	522	LYS

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Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	56	ASN
1	А	193	GLN
1	А	194	ASN

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)

2 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



	Mol Type		Mol	Type Chain Res		o Chain	Res	Link	Bo	ond leng	\mathbf{ths}	B	ond ang	les
	WIOI	Type Chain Kes Lin		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2					
	2	4AM	А	601	-	20,20,20	1.17	2 (10%)	22,28,28	1.44	3 (13%)			
	3	CIT	А	602	-	12,12,12	1.33	2 (16%)	17,17,17	1.46	3 (17%)			

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).

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	4AM	А	601	-	-	3/18/34/34	0/1/1/1
3	CIT	А	602	-	-	6/16/16/16	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	А	602	CIT	C3-C6	2.80	1.56	1.53
2	А	601	4AM	O1B-C1	-2.64	1.22	1.30
2	А	601	4AM	C8-C7	2.33	1.57	1.53
3	А	602	CIT	O4-C5	-2.06	1.23	1.30

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
2	А	601	4AM	C8-C7-C6	4.24	121.07	113.03
3	А	602	CIT	O5-C6-C3	-3.44	117.38	122.25
3	А	602	CIT	O6-C6-C3	2.71	117.76	113.05
2	А	601	4AM	O6-C2-C1	2.51	117.08	112.06
2	А	601	4AM	C6-C5-N5	2.23	114.62	110.91
3	А	602	CIT	O1-C1-C2	-2.20	116.52	122.94

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	А	602	CIT	C4-C3-C6-O5
3	А	602	CIT	C4-C3-C6-O6
2	А	601	4AM	O8-C8-C9-O9

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Mol	Chain	Res	Type	Atoms
2	А	601	4AM	C7-C8-C9-O9
3	А	602	CIT	O7-C3-C6-O6
3	А	602	CIT	C2-C3-C6-O6
3	А	602	CIT	O7-C3-C6-O5
3	А	602	CIT	C2-C3-C6-O5
2	А	601	4AM	C11-C10-N5-C5

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There are no ring outliers.

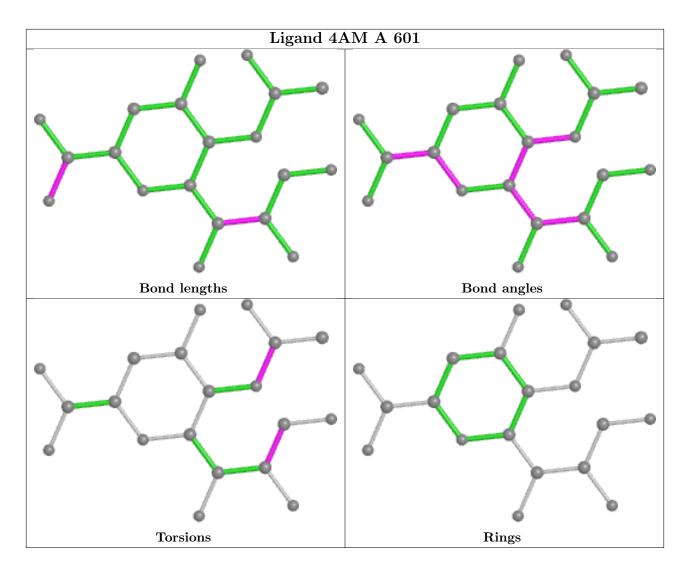
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	А	601	4AM	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 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	Analysed	$\langle RSRZ \rangle$	#RSRZ>2	$OWAB(Å^2)$	Q<0.9
1	А	512/544~(94%)	-0.08	13 (2%) 57 59	30, 48, 84, 145	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	359	PHE	5.1
1	А	79	TYR	3.5
1	А	309	ASN	3.4
1	А	306	LYS	3.2
1	А	307	LYS	3.1
1	А	76	LEU	3.1
1	А	426	ASN	2.8
1	А	77	LYS	2.7
1	А	80	ILE	2.6
1	А	63	ILE	2.2
1	А	83	LYS	2.2
1	А	78	PRO	2.1
1	А	387	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 monosaccharides in this entry.

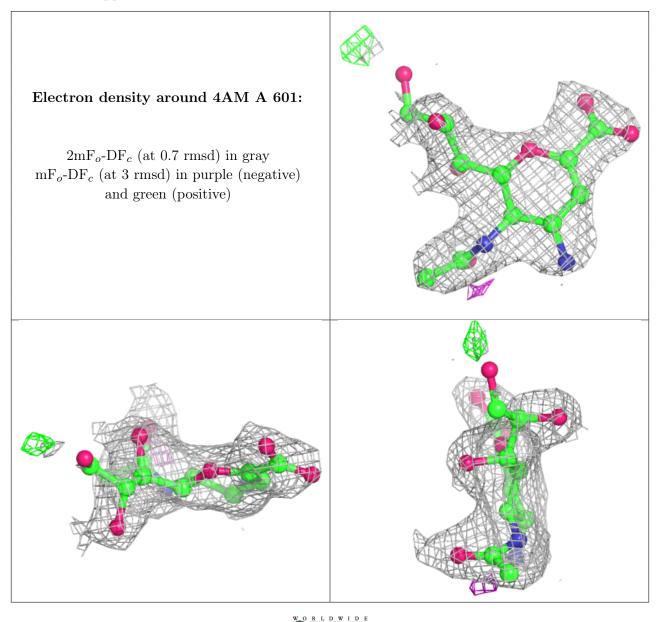


6.4 Ligands (i)

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	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q<0.9
2	4AM	А	601	20/20	0.84	0.21	58,88,95,96	0
3	CIT	А	602	13/13	0.86	0.18	61,81,91,93	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.

