



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 21, 2023 – 05:22 AM JST

PDB ID : 7ESV
Title : Structure and mutation analysis of the hexameric P4 from Pseudomonas aeruginosa phage phiYY
Authors : Zhang, C.Y.; Jin, T.C.
Deposited on : 2021-05-11
Resolution : 2.18 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

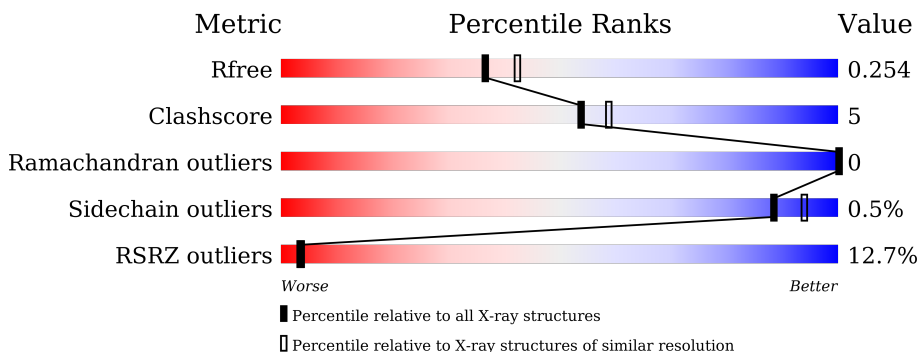
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.18 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	6864 (2.20-2.16)
Clashscore	141614	7689 (2.20-2.16)
Ramachandran outliers	138981	7564 (2.20-2.16)
Sidechain outliers	138945	7564 (2.20-2.16)
RSRZ outliers	127900	6738 (2.20-2.16)

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	291	 12% 80% 8% 11%
1	B	291	 10% 78% 12% 9%
1	C	291	 12% 78% 14% 8%

2 Entry composition [i](#)

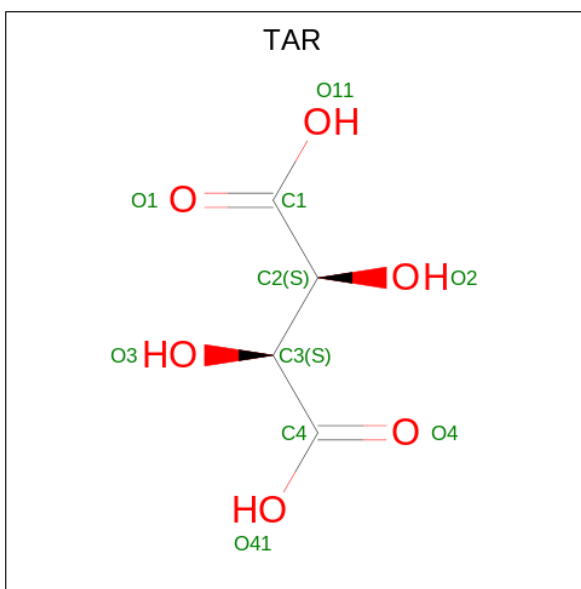
There are 3 unique types of molecules in this entry. The entry contains 5915 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Packaging NTPase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	B	264	Total 1944	C 1226	N 335	O 371	S 12	0	0	0
1	A	258	Total 1909	C 1207	N 326	O 364	S 12	0	0	0
1	C	267	Total 1986	C 1253	N 345	O 376	S 12	0	1	0

- Molecule 2 is D(-)-TARTARIC ACID (three-letter code: TAR) (formula: C₄H₆O₆) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	B	1	Total 10	C 4	O 6	0	0

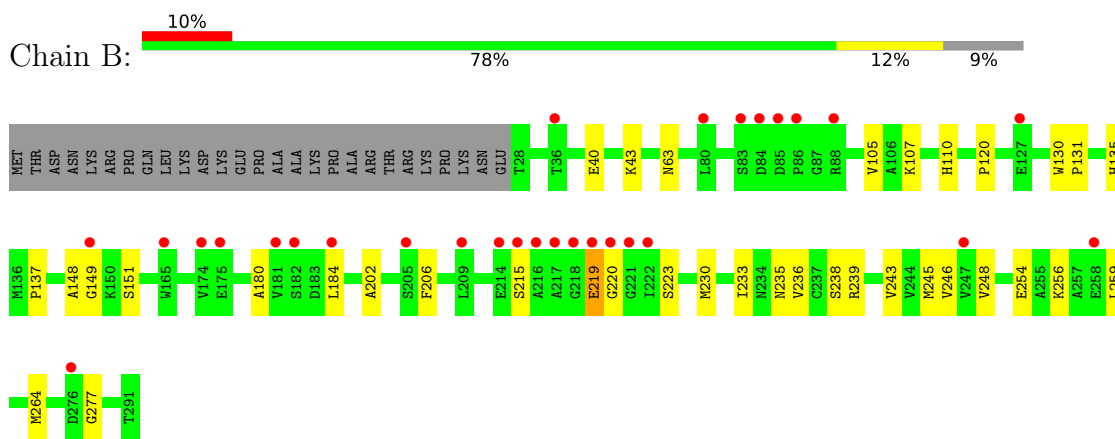
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	23	Total 23	O 23	0	0
3	A	22	Total 22	O 22	0	0
3	C	21	Total 21	O 21	0	0

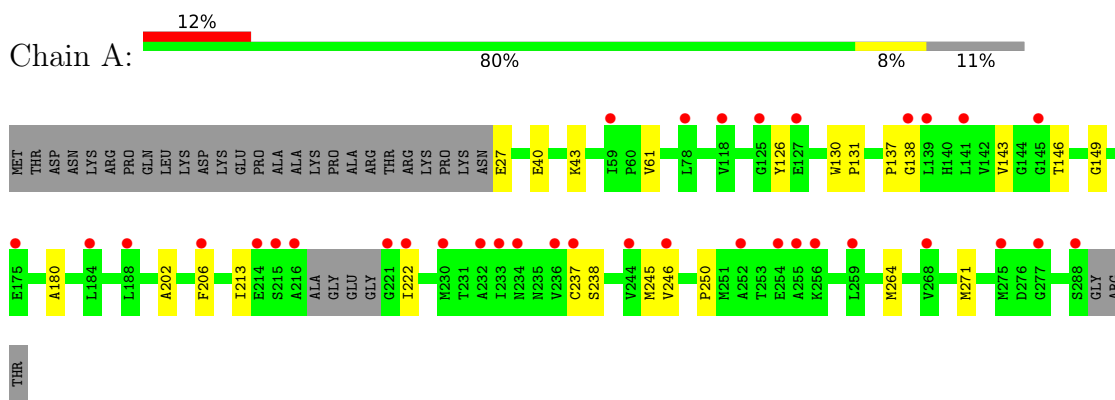
3 Residue-property plots

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.

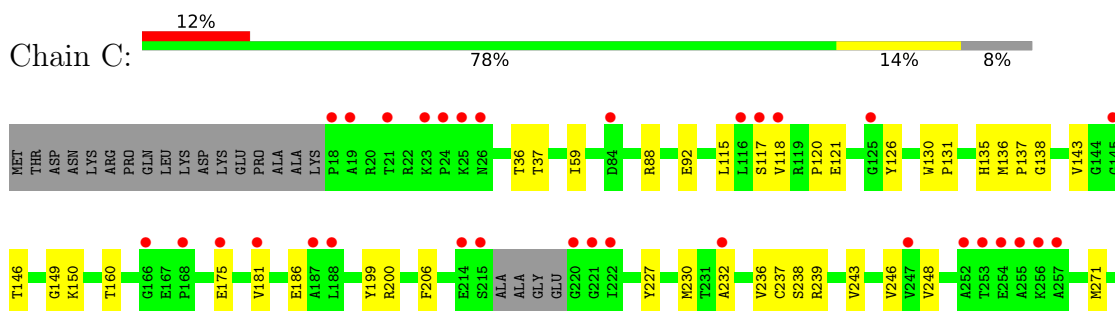
- Molecule 1: Packaging NTPase



- Molecule 1: Packaging NTPase



- Molecule 1: Packaging NTPase



L274	R275	D276	G277	A278	V279	S288	GLY	ARG	THR
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4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	155.63Å 67.93Å 107.65Å 90.00° 128.94° 90.00°	Depositor
Resolution (Å)	33.57 – 2.18 33.57 – 2.18	Depositor EDS
% Data completeness (in resolution range)	99.3 (33.57-2.18) 99.3 (33.57-2.18)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.21 (at 2.18Å)	Xtrriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.212 , 0.255 0.212 , 0.254	Depositor DCC
R_{free} test set	1988 reflections (4.37%)	wwPDB-VP
Wilson B-factor (Å ²)	42.7	Xtrriage
Anisotropy	0.733	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 37.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5915	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.77% 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: TAR

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	0.28	0/1943	0.52	0/2648
1	B	0.31	0/1979	0.53	0/2697
1	C	0.34	0/2022	0.55	0/2752
All	All	0.31	0/5944	0.53	0/8097

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 [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	1909	0	1919	14	0
1	B	1944	0	1954	26	0
1	C	1986	0	2004	25	0
2	B	10	0	4	3	0
3	A	22	0	0	0	0
3	B	23	0	0	0	0
3	C	21	0	0	0	0
All	All	5915	0	5881	62	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 5.

All (62) 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:B:220:GLY:HA2	1:B:256:LYS:HB3	1.60	0.83
1:C:137:PRO:HB2	1:C:238:SER:HA	1.71	0.73
1:A:137:PRO:HB2	1:A:238:SER:HA	1.71	0.72
1:B:215:SER:HB3	1:B:223:SER:HB3	1.76	0.67
1:B:151:SER:H	2:B:301:TAR:H3	1.59	0.67
1:C:88:ARG:O	1:C:92:GLU:HG3	1.96	0.66
1:B:219:GLU:HB3	1:B:259:LEU:HD21	1.78	0.64
1:B:219:GLU:O	1:B:256:LYS:HD3	2.00	0.61
1:C:146:THR:HG22	1:C:149:GLY:H	1.67	0.60
1:C:181:VAL:HB	1:C:186:GLU:HG3	1.87	0.57
1:A:143:VAL:HG21	1:A:271:MET:HE2	1.89	0.55
1:C:150:LYS:HG2	1:C:274:LEU:HD22	1.87	0.55
1:C:160:THR:HG21	1:C:199:TYR:HD1	1.73	0.54
1:B:219:GLU:HB3	1:B:259:LEU:CD2	2.38	0.54
1:A:222:ILE:HD11	1:A:250:PRO:HB2	1.90	0.53
1:C:131:PRO:HD2	1:C:135:HIS:CD2	2.45	0.52
1:A:146:THR:HG23	1:A:149:GLY:H	1.75	0.52
1:B:137:PRO:HB2	1:B:238:SER:HA	1.90	0.52
1:C:232:ALA:O	1:C:236:VAL:HG23	2.11	0.51
1:C:121:GLU:HB2	1:C:239:ARG:HG2	1.92	0.51
1:C:136:MET:HE3	1:C:137:PRO:HD2	1.92	0.51
1:B:130:TRP:CD1	1:B:131:PRO:HA	2.47	0.49
1:C:206:PHE:CD2	1:C:246:VAL:HG11	2.47	0.49
1:A:130:TRP:CD1	1:A:131:PRO:HA	2.49	0.48
1:B:131:PRO:HD2	1:B:135:HIS:CD2	2.49	0.48
1:B:131:PRO:HD3	1:B:243:VAL:HG11	1.96	0.48
1:C:130:TRP:CD2	1:C:200:ARG:HD3	2.49	0.48
1:A:126:TYR:CE1	1:A:137:PRO:HD3	2.49	0.47
1:C:138:GLY:HA2	1:C:237:CYS:HB2	1.96	0.47
1:B:105:VAL:O	1:B:107:LYS:HD2	2.15	0.47
1:B:149:GLY:HA2	2:B:301:TAR:H2	1.97	0.46
1:B:206:PHE:CD2	1:B:246:VAL:HG11	2.50	0.46
1:C:131:PRO:HD3	1:C:243:VAL:HG11	1.96	0.46
1:B:63:ASN:O	1:B:110:HIS:NE2	2.45	0.45
1:C:130:TRP:CD1	1:C:131:PRO:HA	2.52	0.45
1:C:126:TYR:CE2	1:C:137:PRO:HD3	2.53	0.44
1:B:180:ALA:HB1	1:C:120:PRO:HD2	1.99	0.44
1:C:36:THR:HG23	1:C:37:THR:HG23	2.00	0.44
1:B:235:ASN:O	1:B:239:ARG:HG3	2.18	0.43
1:B:149:GLY:CA	2:B:301:TAR:H2	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:184:LEU:HD22	1:C:115:LEU:HD12	2.00	0.43
1:B:148:ALA:O	1:B:277:GLY:N	2.49	0.43
1:C:143:VAL:HG21	1:C:271:MET:HE3	1.99	0.43
1:B:202:ALA:HA	1:B:245:MET:O	2.19	0.43
1:A:61:VAL:HG12	1:A:213:ILE:HD11	2.01	0.43
1:A:27:GLU:N	1:A:27:GLU:OE1	2.52	0.42
1:A:206:PHE:CD2	1:A:246:VAL:HG11	2.54	0.42
1:B:233:ILE:O	1:B:236:VAL:HG12	2.19	0.42
1:B:40:GLU:HA	1:B:43:LYS:HE2	2.02	0.42
1:A:138:GLY:HA2	1:A:237:CYS:HB2	2.02	0.42
1:C:117:SER:OG	1:C:118:VAL:N	2.51	0.42
1:A:202:ALA:HA	1:A:245:MET:O	2.20	0.42
1:A:250:PRO:HG3	1:A:264:MET:HE1	2.01	0.42
1:A:40:GLU:HA	1:A:43:LYS:HE2	2.02	0.41
1:C:206:PHE:HB2	1:C:248:VAL:HG22	2.02	0.41
1:B:230:MET:CE	1:B:264:MET:HG2	2.50	0.41
1:B:230:MET:HE3	1:B:230:MET:HB2	1.92	0.41
1:B:206:PHE:HB2	1:B:248:VAL:HG22	2.02	0.41
1:C:59:ILE:HG12	1:C:236:VAL:HG21	2.03	0.41
1:C:227:TYR:HA	1:C:230:MET:HE3	2.01	0.41
1:B:120:PRO:HD2	1:A:180:ALA:HB1	2.02	0.40
1:C:160:THR:HG21	1:C:199:TYR:CD1	2.53	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	254/291 (87%)	244 (96%)	10 (4%)	0	100	100
1	B	262/291 (90%)	254 (97%)	8 (3%)	0	100	100
1	C	264/291 (91%)	259 (98%)	5 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	780/873 (89%)	757 (97%)	23 (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	201/227 (88%)	201 (100%)	0	100	100
1	B	203/227 (89%)	201 (99%)	2 (1%)	76	85
1	C	209/227 (92%)	208 (100%)	1 (0%)	88	94
All	All	613/681 (90%)	610 (100%)	3 (0%)	88	94

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

Mol	Chain	Res	Type
1	B	219	GLU
1	B	254	GLU
1	C	175	GLU

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](#)

1 ligand is 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
2	TAR	B	301	-	9,9,9	0.99	0	12,12,12	1.28	2 (16%)

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	TAR	B	301	-	-	4/12/12/12	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	TAR	O41-C4-C3	2.46	119.92	113.27
2	B	301	TAR	O11-C1-C2	2.40	119.76	113.27

There are no chirality outliers.

All (4) torsion outliers are listed below:

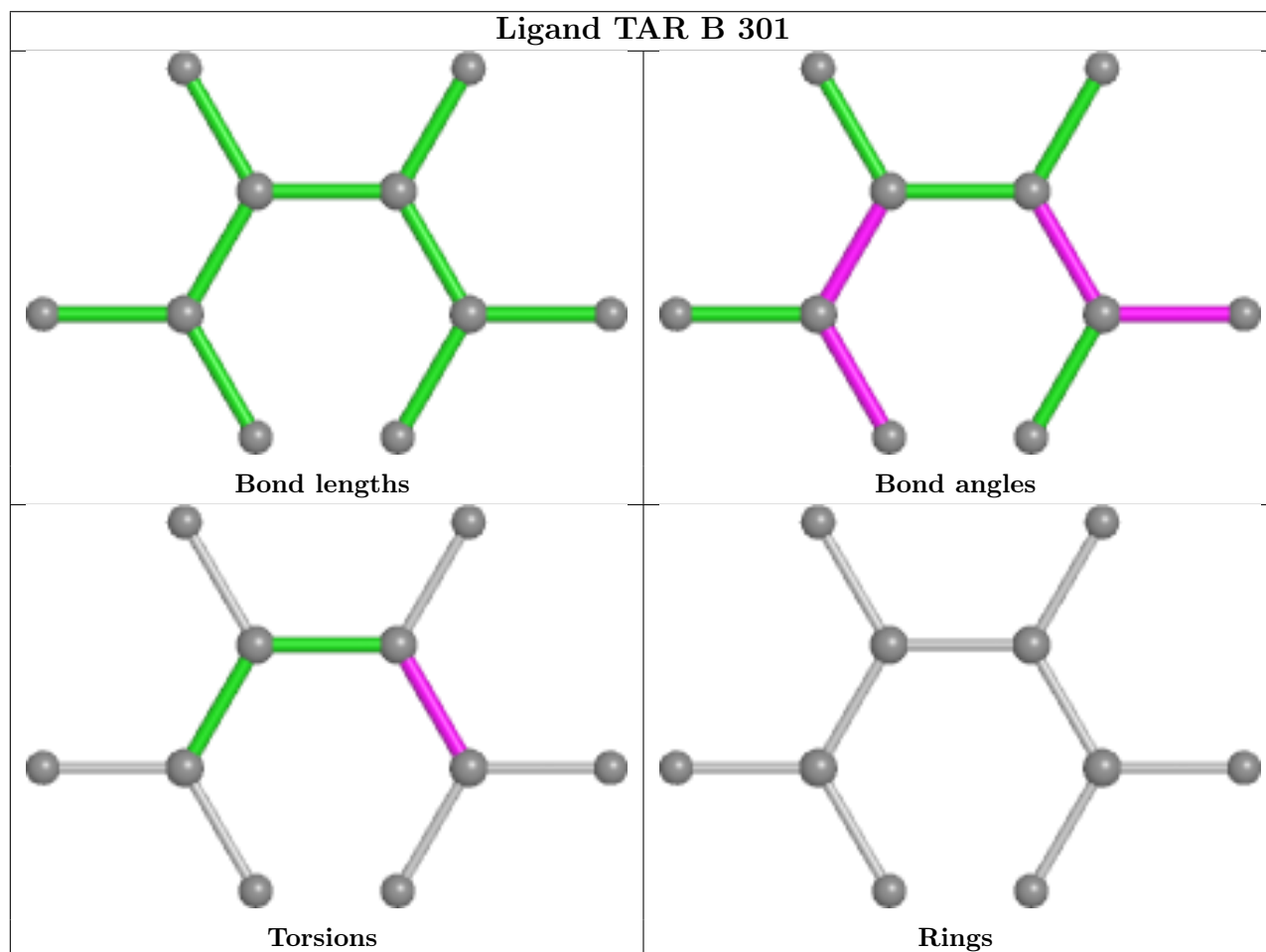
Mol	Chain	Res	Type	Atoms
2	B	301	TAR	O1-C1-C2-C3
2	B	301	TAR	O11-C1-C2-C3
2	B	301	TAR	O1-C1-C2-O2
2	B	301	TAR	O11-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	301	TAR	3	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, 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	258/291 (88%)	0.66	35 (13%) 3 3	38, 47, 74, 82	0
1	B	264/291 (90%)	0.66	29 (10%) 5 6	36, 47, 72, 92	0
1	C	267/291 (91%)	0.80	36 (13%) 3 3	38, 51, 79, 95	0
All	All	789/873 (90%)	0.71	100 (12%) 3 3	36, 48, 75, 95	0

All (100) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	220	GLY	8.3
1	B	217	ALA	7.5
1	B	222	ILE	6.6
1	A	222	ILE	6.1
1	C	25	LYS	6.0
1	C	24	PRO	5.8
1	C	18	PRO	5.4
1	A	233	ILE	5.0
1	B	216	ALA	4.9
1	C	19	ALA	4.8
1	B	219	GLU	4.8
1	B	215	SER	4.7
1	C	215	SER	4.6
1	A	215	SER	4.5
1	C	222	ILE	4.4
1	C	276	ASP	4.3
1	C	145	GLY	4.2
1	A	216	ALA	3.9
1	B	165	TRP	3.9
1	A	125	GLY	3.9
1	B	214	GLU	3.9
1	B	218	GLY	3.9
1	A	221	GLY	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	139	LEU	3.9
1	B	84	ASP	3.8
1	A	275	MET	3.7
1	B	221	GLY	3.7
1	C	21	THR	3.6
1	C	221	GLY	3.4
1	A	255	ALA	3.4
1	A	175	GLU	3.4
1	A	252	ALA	3.3
1	A	237	CYS	3.3
1	A	256	LYS	3.2
1	C	254	GLU	3.2
1	B	174	VAL	3.2
1	C	255	ALA	3.1
1	C	125	GLY	3.1
1	B	83	SER	3.0
1	B	205	SER	3.0
1	A	188	LEU	3.0
1	B	220	GLY	3.0
1	C	252	ALA	3.0
1	C	23	LYS	2.9
1	B	149	GLY	2.9
1	C	181	VAL	2.8
1	C	118	VAL	2.8
1	C	232	ALA	2.8
1	C	188	LEU	2.8
1	B	88	ARG	2.8
1	C	256	LYS	2.8
1	A	138	GLY	2.8
1	A	184	LEU	2.8
1	A	214	GLU	2.7
1	B	276	ASP	2.7
1	B	247	VAL	2.7
1	C	84	ASP	2.6
1	C	278	ALA	2.6
1	C	166	GLY	2.6
1	A	118	VAL	2.6
1	C	117	SER	2.6
1	A	288	SER	2.5
1	A	277	GLY	2.5
1	C	253	THR	2.5
1	A	145	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	86	PRO	2.5
1	A	244	VAL	2.5
1	A	206	PHE	2.5
1	C	275	MET	2.4
1	B	258	GLU	2.4
1	C	257	ALA	2.4
1	A	254	GLU	2.4
1	A	259	LEU	2.4
1	B	85	ASP	2.4
1	A	127	GLU	2.4
1	A	232	ALA	2.4
1	B	80	LEU	2.4
1	A	78	LEU	2.4
1	A	246	VAL	2.4
1	A	268	VAL	2.3
1	A	234	ASN	2.3
1	B	184	LEU	2.3
1	C	214	GLU	2.3
1	C	247	VAL	2.3
1	C	175	GLU	2.3
1	A	236	VAL	2.3
1	C	187	ALA	2.2
1	C	279	VAL	2.2
1	A	230	MET	2.2
1	B	175	GLU	2.1
1	B	209	LEU	2.1
1	B	182	SER	2.0
1	A	59	ILE	2.0
1	A	141	LEU	2.0
1	C	26	ASN	2.0
1	B	127	GLU	2.0
1	B	181	VAL	2.0
1	C	168	PRO	2.0
1	B	36	THR	2.0
1	C	116	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates [i](#)

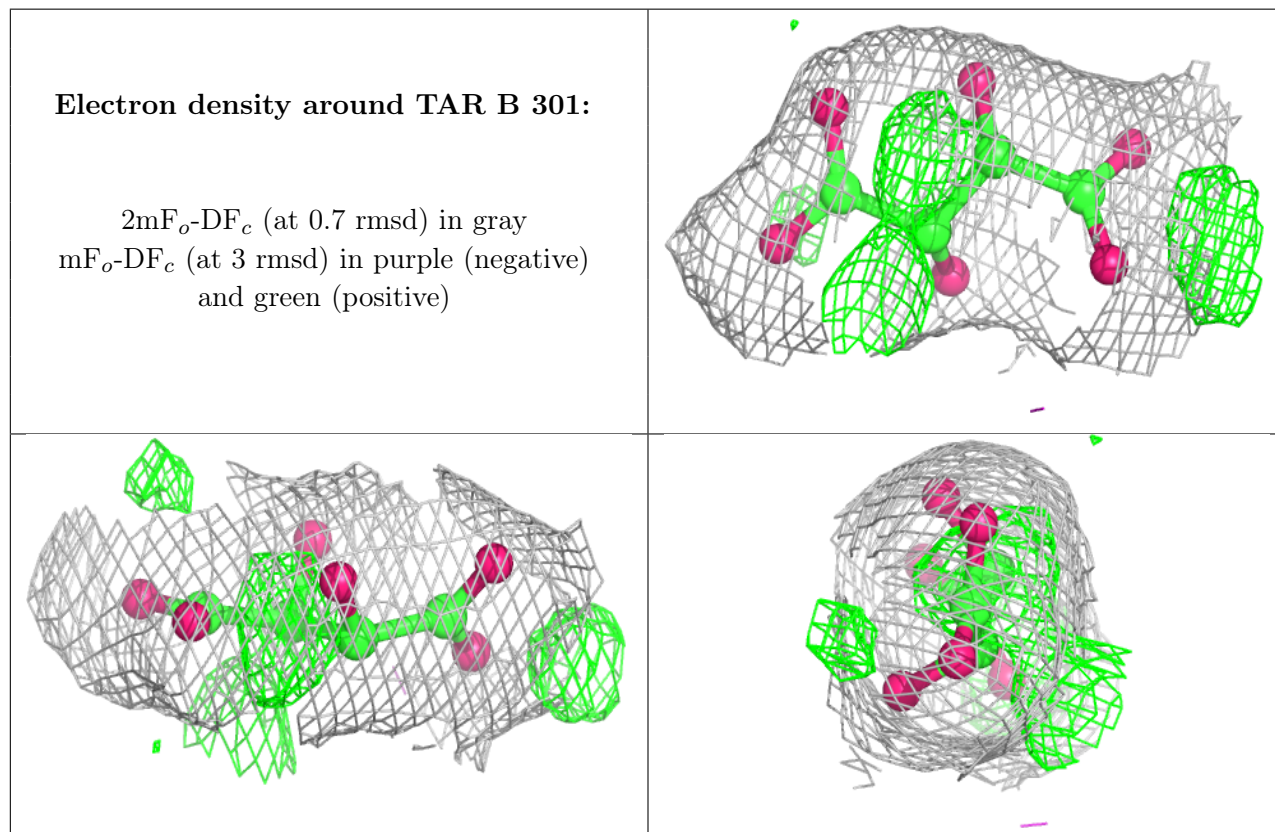
There are no monosaccharides in this entry.

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
2	TAR	B	301	10/10	0.90	0.21	61,64,72,73	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.