



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 2, 2026 – 12:50 AM UTC

PDB ID : 9RJK / pdb_00009rjk
Title : Structure of Mycobacterium tuberculosis InhA in complex with pyridomycin derivative KV41a (compound 11)
Authors : Publicola, G.; Mourey, L.; Valderrama, K.; Hartkoorn, R.; Maveyraud, L.
Deposited on : 2025-06-12
Resolution : 1.66 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

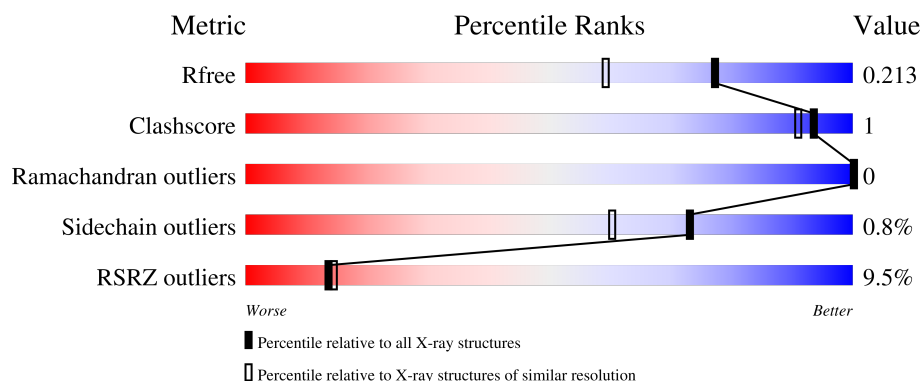
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.66 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}	180053	2563 (1.66-1.66)
Clashscore	190562	2662 (1.66-1.66)
Ramachandran outliers	187476	2621 (1.66-1.66)
Sidechain outliers	187428	2621 (1.66-1.66)
RSRZ outliers	180081	2564 (1.66-1.66)

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	272	5% 89% 7%
1	B	272	5% 90% 7%
1	C	272	12% 92%
1	D	272	13% 89% 8%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8671 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.

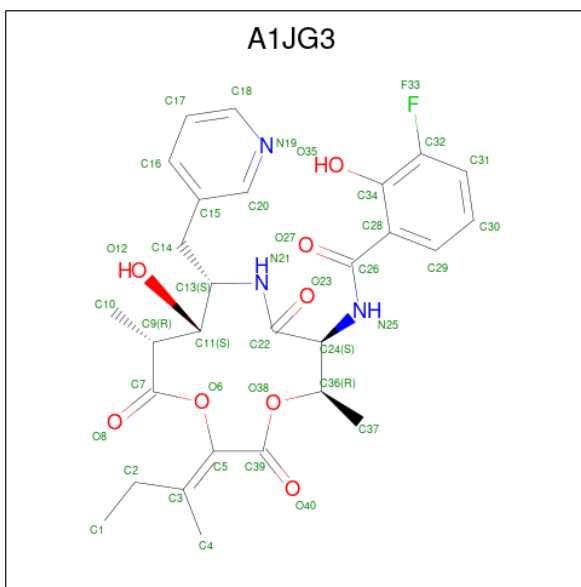
- Molecule 1 is a protein called Enoyl-[acyl-carrier-protein] reductase [NADH].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	254	Total	C	N	O	S	0	3	0
			1927	1220	340	357	10			
1	B	252	Total	C	N	O	S	0	5	0
			1925	1218	339	357	11			
1	C	263	Total	C	N	O	S	0	3	0
			1968	1240	347	371	10			
1	D	251	Total	C	N	O	S	0	3	0
			1861	1180	323	348	10			

There are 12 discrepancies between the modelled and reference sequences:

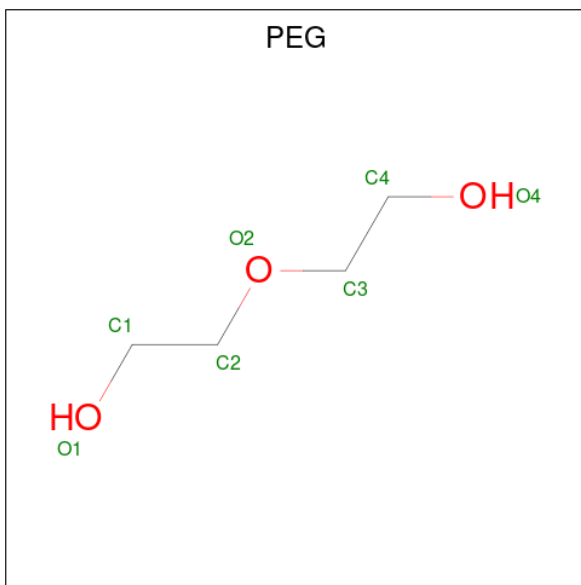
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P9WGR1
A	-1	SER	-	expression tag	UNP P9WGR1
A	0	HIS	-	expression tag	UNP P9WGR1
B	-2	GLY	-	expression tag	UNP P9WGR1
B	-1	SER	-	expression tag	UNP P9WGR1
B	0	HIS	-	expression tag	UNP P9WGR1
C	-2	GLY	-	expression tag	UNP P9WGR1
C	-1	SER	-	expression tag	UNP P9WGR1
C	0	HIS	-	expression tag	UNP P9WGR1
D	-2	GLY	-	expression tag	UNP P9WGR1
D	-1	SER	-	expression tag	UNP P9WGR1
D	0	HIS	-	expression tag	UNP P9WGR1

- Molecule 2 is {N}-(2 {Z},5 {R},6 {S},9 {S},10 {S},11 {R})-2-butan-2-ylidene-5,11-dimethyl-10-oxidanyl-3,7,12-tris(oxidanylidene)-9-(pyridin-3-ylmethyl)-1,4-dioxo-8-azacyclododec-6-yl]-3-fluoranyl-2-oxidanyl-benzamide (CCD ID: A1JG3) (formula: C₂₈H₃₂FN₃O₈) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 40	C 28	F 1	N 3	O 8	0	0
2	B	1	Total 40	C 28	F 1	N 3	O 8	0	0
2	C	1	Total 40	C 28	F 1	N 3	O 8	0	0
2	D	1	Total 40	C 28	F 1	N 3	O 8	0	0

- Molecule 3 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			7	4	3		

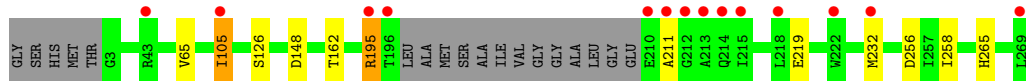
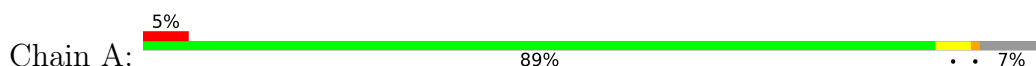
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	246	Total	O	0	0
			246	246		
4	B	239	Total	O	0	0
			239	239		
4	C	181	Total	O	0	0
			181	181		
4	D	150	Total	O	0	0
			150	150		

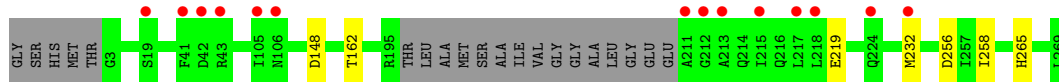
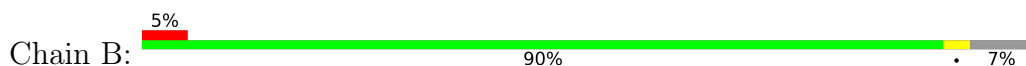
3 Residue-property plots [i](#)

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.

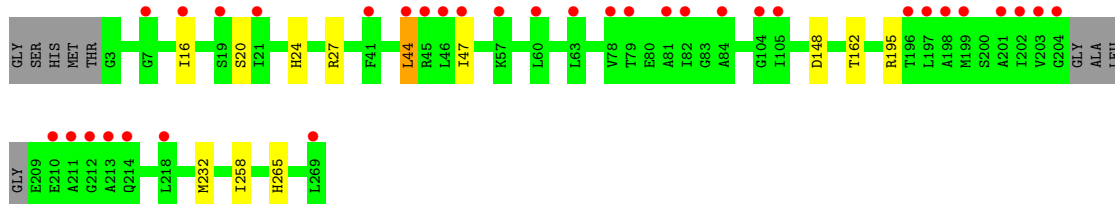
- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]



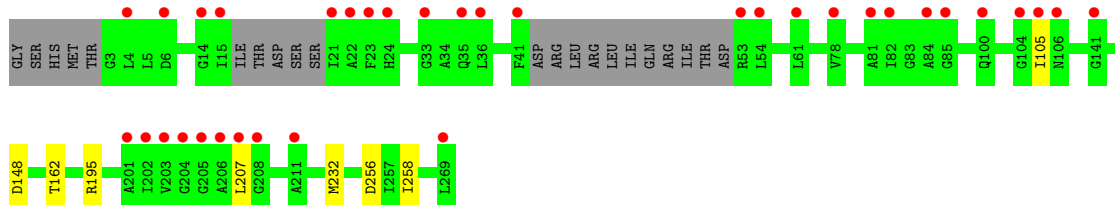
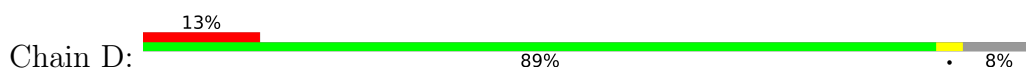
- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]



- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]



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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	82.07Å 81.35Å 95.48Å 90.00° 99.41° 90.00°	Depositor
Resolution (Å)	94.19 – 1.66 94.19 – 1.66	Depositor EDS
% Data completeness (in resolution range)	99.7 (94.19-1.66) 99.6 (94.19-1.66)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 1.66Å)	Xtriage
Refinement program	BUSTER 2.10.4	Depositor
R, R_{free}	0.185 , 0.205 (Not available) , 0.213	Depositor DCC
R_{free} test set	7362 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	25.9	Xtriage
Anisotropy	0.440	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 52.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	8671	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.72% 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: A1JG3, PEG

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	0.82	0/1964	0.95	3/2663 (0.1%)
1	B	0.82	0/1962	0.95	1/2662 (0.0%)
1	C	0.74	0/2004	0.96	1/2721 (0.0%)
1	D	0.70	0/1896	0.92	1/2573 (0.0%)
All	All	0.77	0/7826	0.95	6/10619 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	148	ASP	CA-CB-CG	6.78	119.38	112.60
1	D	148	ASP	CA-CB-CG	6.33	118.93	112.60
1	C	148	ASP	CA-CB-CG	5.89	118.49	112.60
1	A	148	ASP	CA-CB-CG	5.50	118.10	112.60
1	A	195	ARG	CA-C-N	5.41	131.43	121.70
1	A	195	ARG	C-N-CA	5.41	131.43	121.70

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	1927	0	1930	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1925	0	1923	5	0
1	C	1968	0	1952	8	0
1	D	1861	0	1846	4	0
2	A	40	0	0	0	0
2	B	40	0	0	0	0
2	C	40	0	0	0	0
2	D	40	0	0	0	0
3	A	7	0	10	0	0
3	B	7	0	10	0	0
4	A	246	0	0	1	0
4	B	239	0	0	0	0
4	C	181	0	0	2	0
4	D	150	0	0	0	0
All	All	8671	0	7671	21	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 1.

All (21) 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:C:24:HIS:HD2	1:C:27:ARG:HE	1.43	0.65
1:C:20:SER:CB	4:C:404:HOH:O	2.46	0.63
1:D:195:ARG:HG2	1:D:232:MET:HE2	1.82	0.61
1:A:219:GLU:O	4:A:401:HOH:O	2.16	0.58
1:D:105:ILE:CB	1:D:207:LEU:CB	2.82	0.57
1:A:195:ARG:HB2	1:A:232:MET:HE2	1.90	0.54
1:C:265:HIS:HE1	1:D:256:ASP:OD2	1.91	0.54
1:C:16:ILE:HG12	4:C:404:HOH:O	2.08	0.53
1:A:219:GLU:OE2	1:A:232:MET:HE1	2.10	0.52
1:C:44:LEU:HD12	1:C:47:ILE:HD12	1.91	0.51
1:A:256:ASP:OD2	1:B:265:HIS:HE1	1.94	0.50
1:B:219:GLU:HB3	1:B:232[B]:MET:SD	2.52	0.50
1:A:105:ILE:HD13	1:A:211:ALA:HB2	1.94	0.49
1:A:265:HIS:HE1	1:B:256:ASP:OD2	1.97	0.48
1:C:24:HIS:CD2	1:C:27:ARG:HE	2.29	0.46
1:C:195:ARG:HG2	1:C:232:MET:HE2	1.99	0.44
1:B:232[B]:MET:HE2	1:B:232[B]:MET:HB2	1.82	0.43
1:A:258:ILE:HD12	1:B:258:ILE:HD12	2.00	0.42
1:A:65:VAL:HB	1:A:126:SER:HB2	2.03	0.41
1:C:258:ILE:HD12	1:D:258:ILE:HD12	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:ILE:HB	1:A:211:ALA:HB2	2.03	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	253/272 (93%)	244 (96%)	9 (4%)	0	100	100
1	B	253/272 (93%)	244 (96%)	9 (4%)	0	100	100
1	C	262/272 (96%)	254 (97%)	8 (3%)	0	100	100
1	D	248/272 (91%)	239 (96%)	9 (4%)	0	100	100
All	All	1016/1088 (93%)	981 (97%)	35 (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	197/207 (95%)	195 (99%)	2 (1%)	68	52
1	B	197/207 (95%)	196 (100%)	1 (0%)	81	72
1	C	199/207 (96%)	197 (99%)	2 (1%)	68	52
1	D	185/207 (89%)	184 (100%)	1 (0%)	81	72

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	778/828 (94%)	772 (99%)	6 (1%)	73	60

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	105	ILE
1	A	162	THR
1	B	162	THR
1	C	44	LEU
1	C	162	THR
1	D	162	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	HIS
1	A	48	GLN
1	A	66	GLN
1	A	265	HIS
1	B	48	GLN
1	B	66	GLN
1	B	121	HIS
1	B	265	HIS
1	C	24	HIS
1	C	66	GLN
1	C	86	ASN
1	C	121	HIS
1	C	265	HIS
1	D	66	GLN
1	D	121	HIS
1	D	139	ASN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 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 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	A1JG3	D	301	-	40,42,42	0.23	0	49,59,59	0.64	1 (2%)
3	PEG	A	302	-	6,6,6	0.12	0	5,5,5	0.12	0
2	A1JG3	A	301	-	40,42,42	0.27	0	49,59,59	0.67	1 (2%)
3	PEG	B	302	-	6,6,6	0.10	0	5,5,5	0.18	0
2	A1JG3	B	301	-	40,42,42	0.22	0	49,59,59	0.64	0
2	A1JG3	C	301	-	40,42,42	0.28	0	49,59,59	0.61	0

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	A1JG3	D	301	-	-	6/53/54/54	0/2/3/3
3	PEG	A	302	-	-	1/4/4/4	-
2	A1JG3	A	301	-	-	4/53/54/54	0/2/3/3
3	PEG	B	302	-	-	1/4/4/4	-
2	A1JG3	B	301	-	-	5/53/54/54	0/2/3/3
2	A1JG3	C	301	-	-	7/53/54/54	0/2/3/3

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	A1JG3	C22-C24-N25	-2.41	103.77	110.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	301	A1JG3	C22-C24-N25	-2.05	104.75	110.32

There are no chirality outliers.

All (24) torsion outliers are listed below:

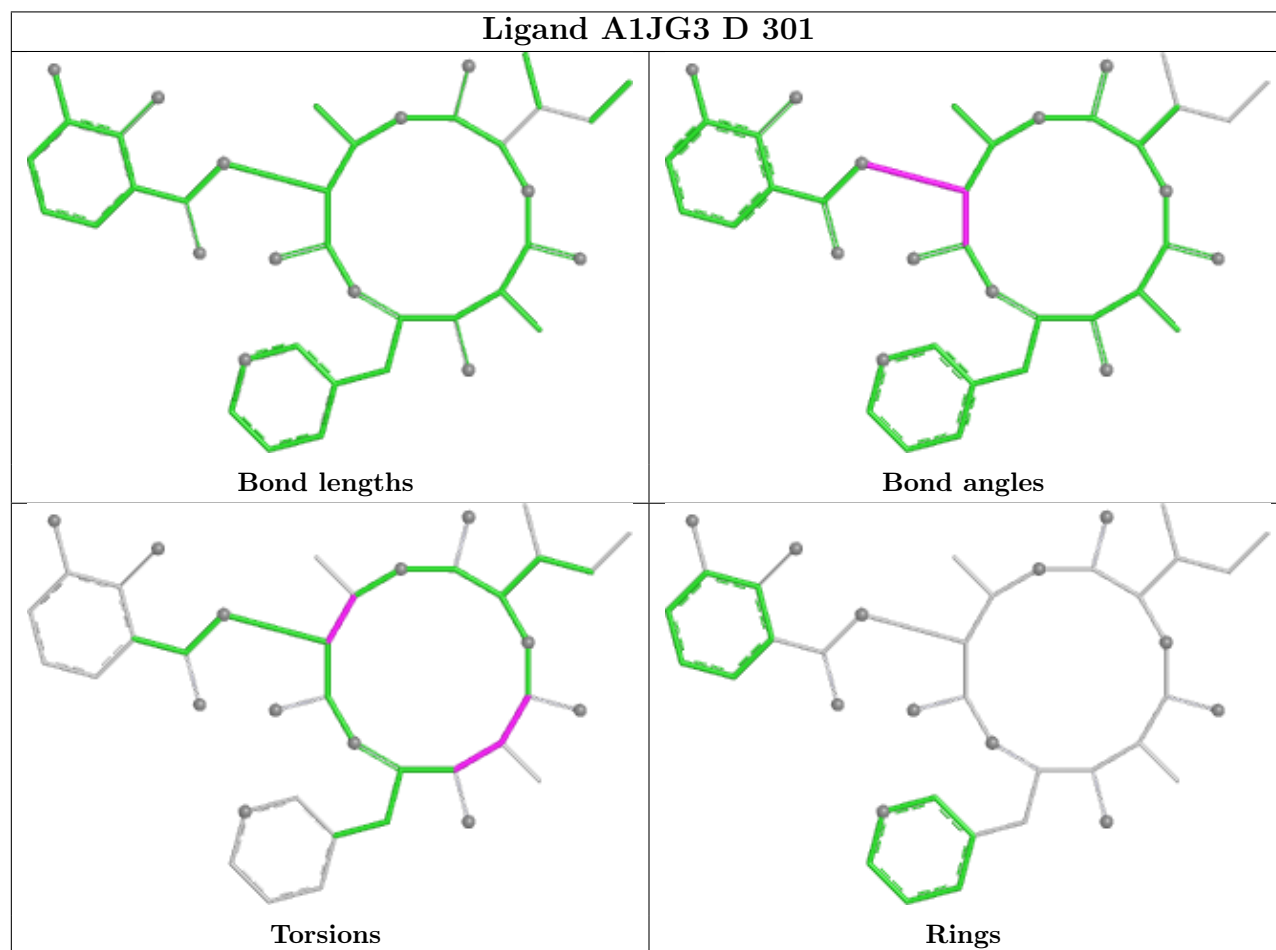
Mol	Chain	Res	Type	Atoms
2	C	301	A1JG3	C22-C24-C36-C37
2	D	301	A1JG3	C22-C24-C36-C37
2	A	301	A1JG3	N25-C24-C36-O38
3	A	302	PEG	C4-C3-O2-C2
3	B	302	PEG	C4-C3-O2-C2
2	A	301	A1JG3	C22-C24-C36-C37
2	B	301	A1JG3	C22-C24-C36-C37
2	A	301	A1JG3	C22-C24-C36-O38
2	B	301	A1JG3	C22-C24-C36-O38
2	C	301	A1JG3	C22-C24-C36-O38
2	D	301	A1JG3	C22-C24-C36-O38
2	C	301	A1JG3	O6-C7-C9-C11
2	D	301	A1JG3	O6-C7-C9-C11
2	A	301	A1JG3	C1-C2-C3-C4
2	C	301	A1JG3	C1-C2-C3-C4
2	B	301	A1JG3	N25-C24-C36-O38
2	D	301	A1JG3	N25-C24-C36-O38
2	C	301	A1JG3	O8-C7-C9-C11
2	D	301	A1JG3	O8-C7-C9-C11
2	C	301	A1JG3	N25-C24-C36-O38
2	B	301	A1JG3	N21-C13-C14-C15
2	B	301	A1JG3	C13-C11-C9-C7
2	C	301	A1JG3	C13-C11-C9-C7
2	D	301	A1JG3	C13-C11-C9-C7

There are no ring outliers.

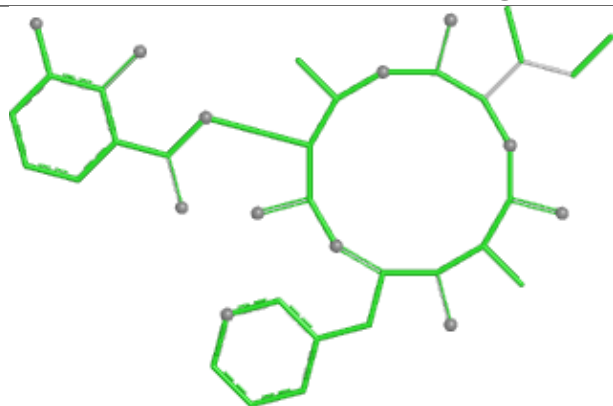
No monomer is involved in short contacts.

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

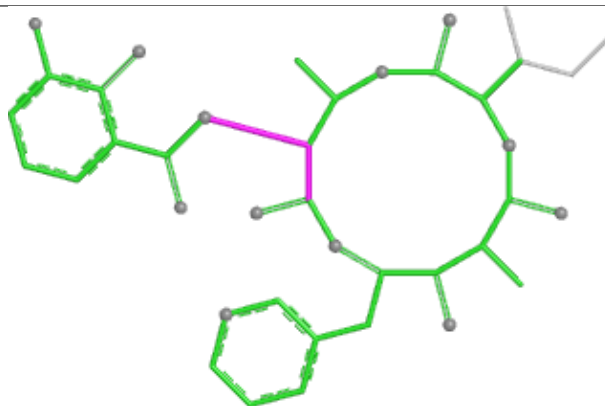
average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



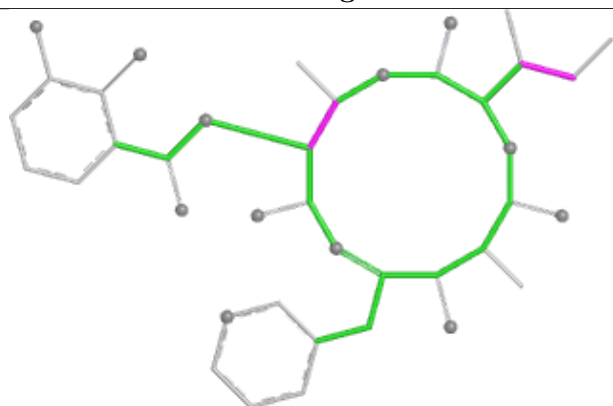
Ligand A1JG3 A 301



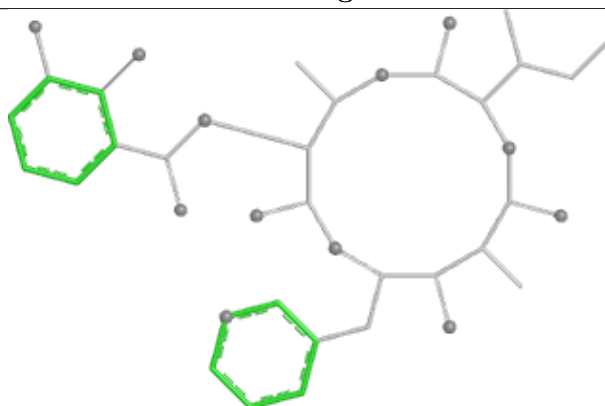
Bond lengths



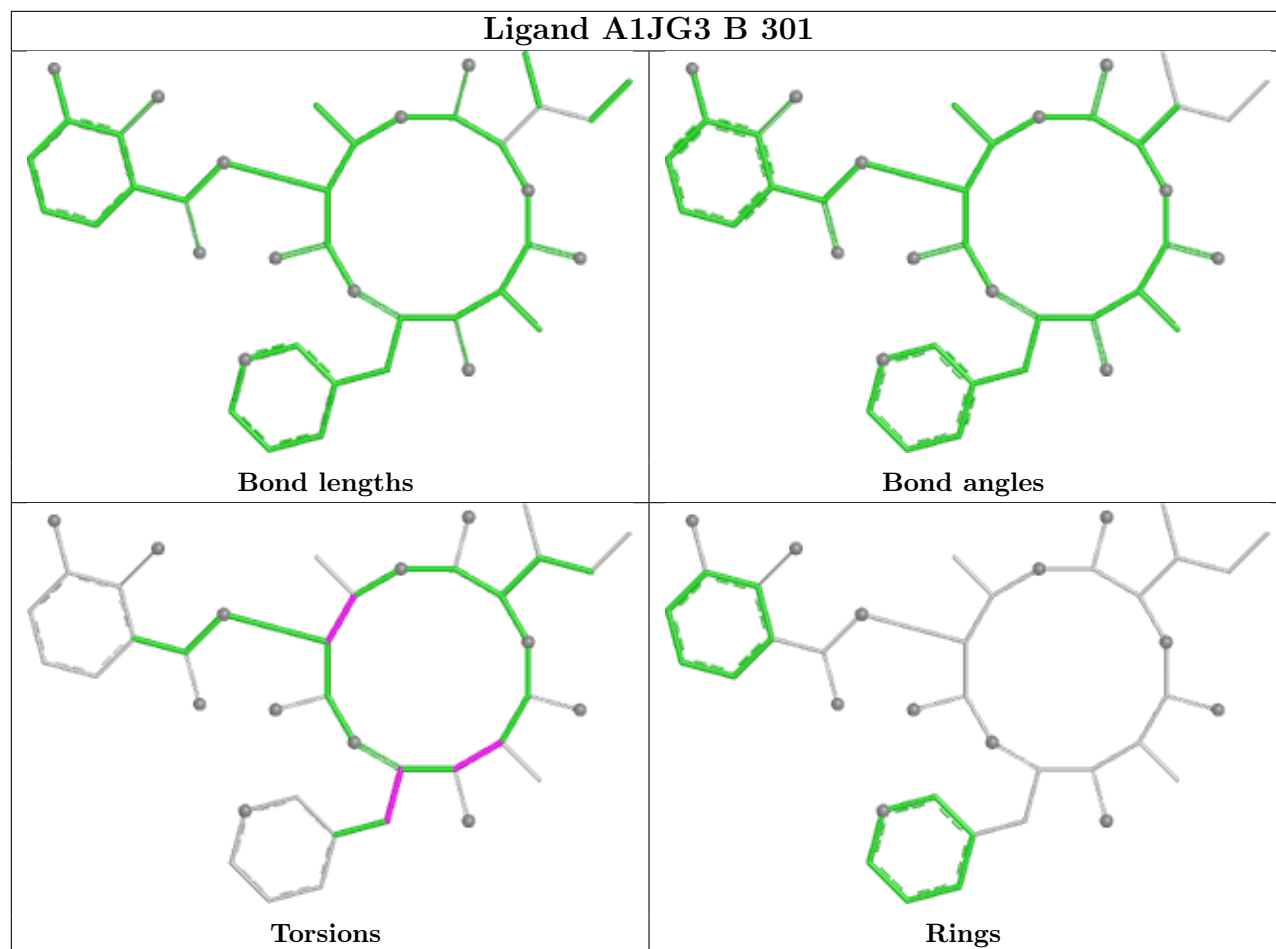
Bond angles

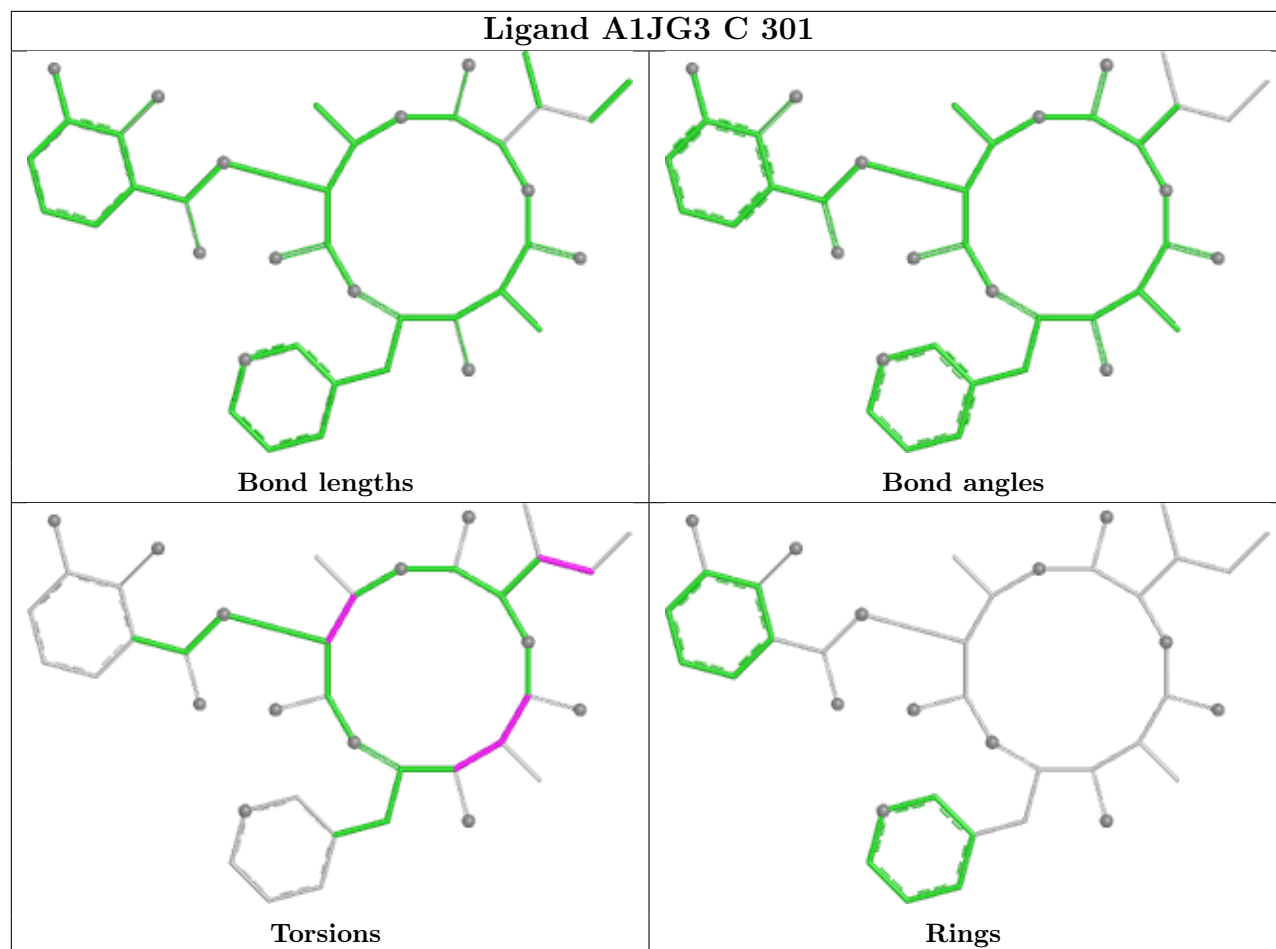


Torsions



Rings





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	254/272 (93%)	0.20	14 (5%) 30 34	12, 31, 50, 75	3 (1%)
1	B	252/272 (92%)	0.14	14 (5%) 30 34	12, 31, 47, 60	5 (1%)
1	C	263/272 (96%)	0.77	34 (12%) 7 8	11, 43, 71, 88	3 (1%)
1	D	251/272 (92%)	0.82	35 (13%) 6 7	11, 43, 77, 88	3 (1%)
All	All	1020/1088 (93%)	0.49	97 (9%) 14 14	11, 36, 68, 88	14 (1%)

All (97) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	211	ALA	7.3
1	D	41	PHE	6.3
1	C	198	ALA	5.8
1	C	105	ILE	5.7
1	C	202	ILE	5.3
1	D	105	ILE	5.0
1	B	211	ALA	5.0
1	D	15	ILE	4.9
1	C	41	PHE	4.6
1	A	196	THR	4.5
1	A	215	ILE	4.3
1	D	206	ALA	4.3
1	D	21	ILE	4.3
1	B	212	GLY	4.2
1	D	81	ALA	4.1
1	C	203	VAL	4.1
1	D	104	GLY	4.0
1	A	212	GLY	4.0
1	C	199	MET	3.9
1	C	197	LEU	3.8
1	A	213	ALA	3.8

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Mol	Chain	Res	Type	RSRZ
1	B	215	ILE	3.7
1	D	269	LEU	3.7
1	C	204	GLY	3.7
1	C	21	ILE	3.6
1	C	104	GLY	3.6
1	C	269	LEU	3.6
1	D	208	GLY	3.5
1	C	201	ALA	3.5
1	A	218	LEU	3.5
1	C	44	LEU	3.5
1	C	84	ALA	3.2
1	D	84	ALA	3.2
1	D	203	VAL	3.2
1	C	46	LEU	3.1
1	D	202	ILE	3.1
1	D	54	LEU	3.1
1	B	213	ALA	3.0
1	C	82	ILE	3.0
1	A	214	GLN	3.0
1	D	204	GLY	3.0
1	D	53	ARG	2.9
1	C	213	ALA	2.9
1	C	212	GLY	2.9
1	C	60	LEU	2.8
1	C	45	ARG	2.8
1	D	205	GLY	2.8
1	B	105	ILE	2.8
1	B	19	SER	2.7
1	C	211	ALA	2.7
1	B	106[A]	ASN	2.7
1	D	106[A]	ASN	2.7
1	D	23	PHE	2.6
1	D	6	ASP	2.6
1	A	105	ILE	2.6
1	D	78	VAL	2.6
1	D	4	LEU	2.5
1	C	57	LYS	2.5
1	A	232	MET	2.5
1	C	47	ILE	2.5
1	D	85	GLY	2.5
1	C	16	ILE	2.5
1	D	24	HIS	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	201	ALA	2.4
1	D	211	ALA	2.4
1	C	214	GLN	2.4
1	D	33	GLY	2.3
1	D	141	GLY	2.3
1	D	207	LEU	2.3
1	A	210	GLU	2.3
1	C	7	GLY	2.3
1	B	217	LEU	2.3
1	C	218	LEU	2.3
1	A	43	ARG	2.3
1	A	195	ARG	2.3
1	D	14	GLY	2.3
1	B	232[A]	MET	2.3
1	B	218	LEU	2.2
1	D	61	LEU	2.2
1	B	41	PHE	2.2
1	C	19	SER	2.2
1	B	42	ASP	2.2
1	B	224	GLN	2.2
1	C	79	THR	2.2
1	D	36	LEU	2.2
1	D	22	ALA	2.1
1	A	222	TRP	2.1
1	C	63	LEU	2.1
1	D	82	ILE	2.1
1	C	196	THR	2.1
1	C	81	ALA	2.1
1	C	210	GLU	2.1
1	D	35	GLN	2.0
1	B	43	ARG	2.0
1	A	269	LEU	2.0
1	D	100	GLN	2.0
1	C	78	VAL	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](#)

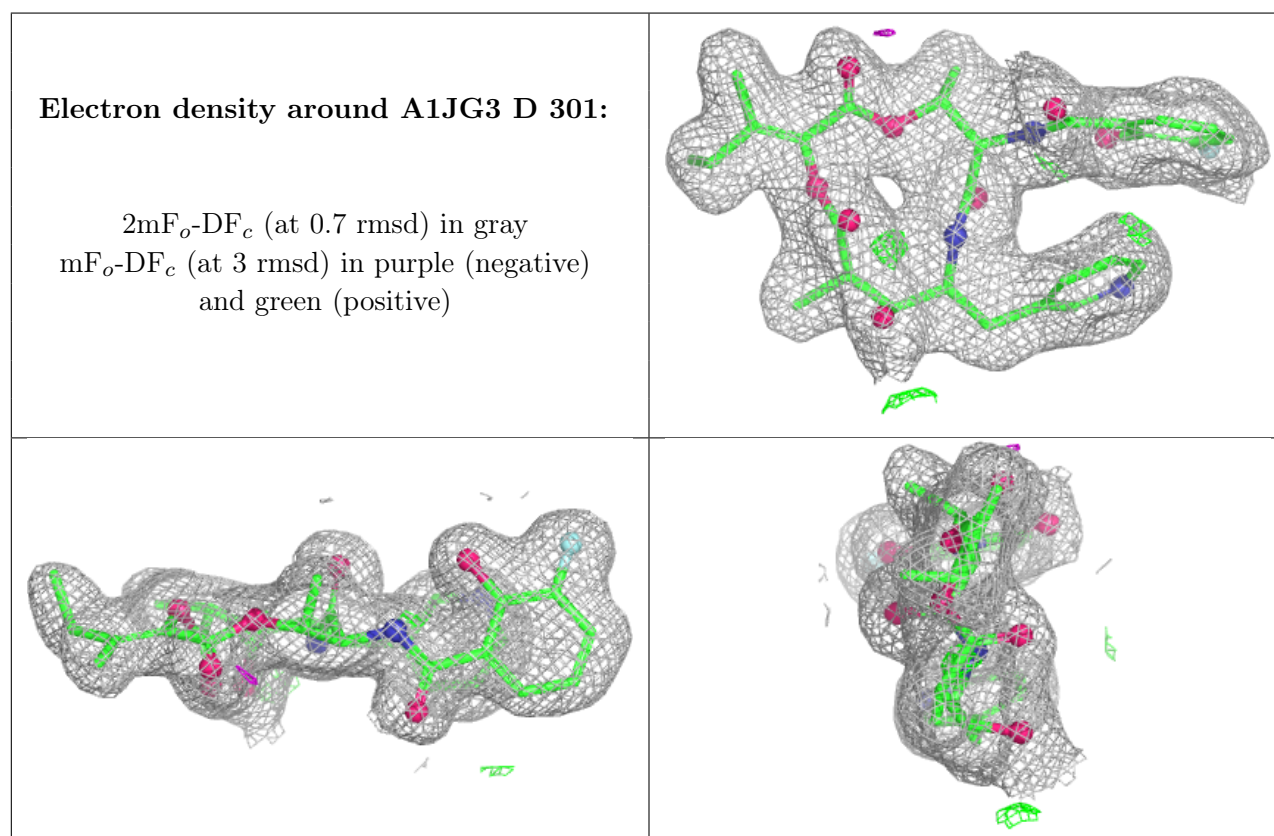
There are no oligosaccharides in this entry.

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.

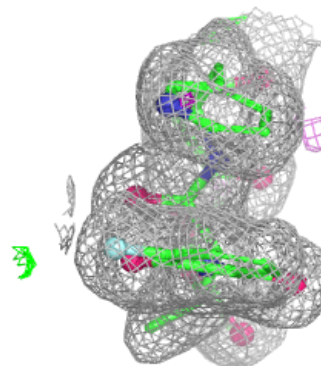
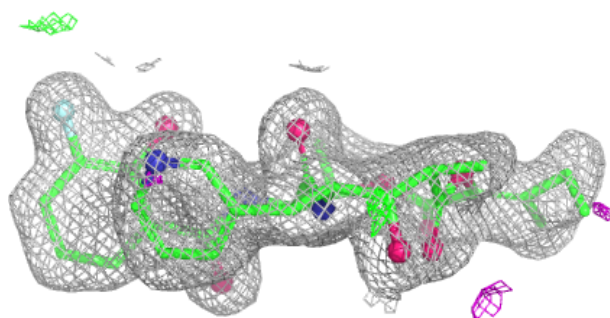
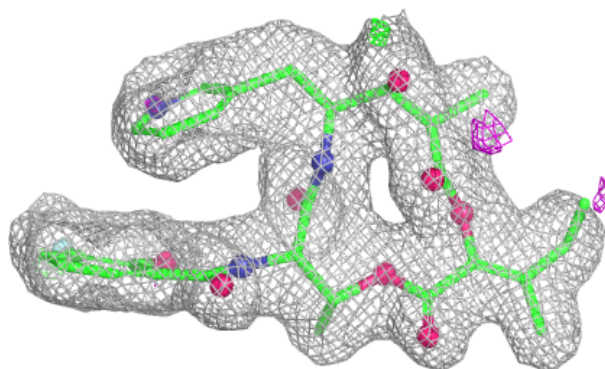
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	PEG	A	302	7/7	0.87	0.14	56,57,59,60	0
3	PEG	B	302	7/7	0.87	0.15	52,54,56,57	0
2	A1JG3	D	301	40/40	0.95	0.08	31,35,37,39	0
2	A1JG3	C	301	40/40	0.96	0.07	30,33,38,41	0
2	A1JG3	B	301	40/40	0.96	0.07	24,27,35,37	0
2	A1JG3	A	301	40/40	0.97	0.06	25,28,34,35	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.

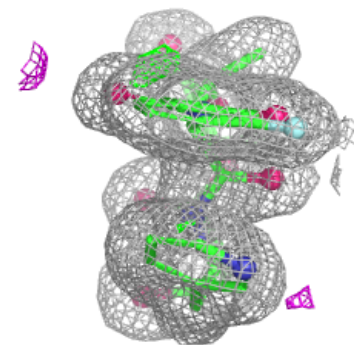
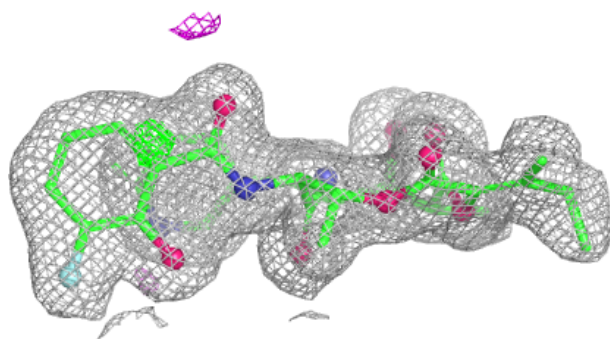
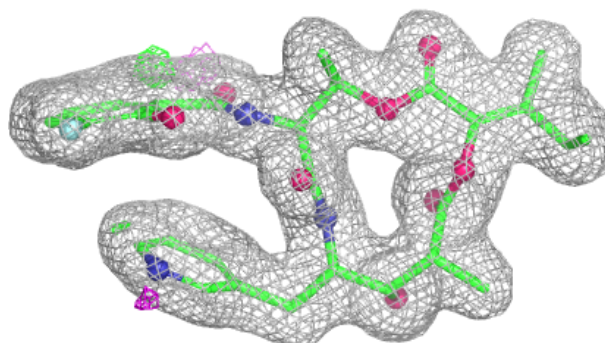


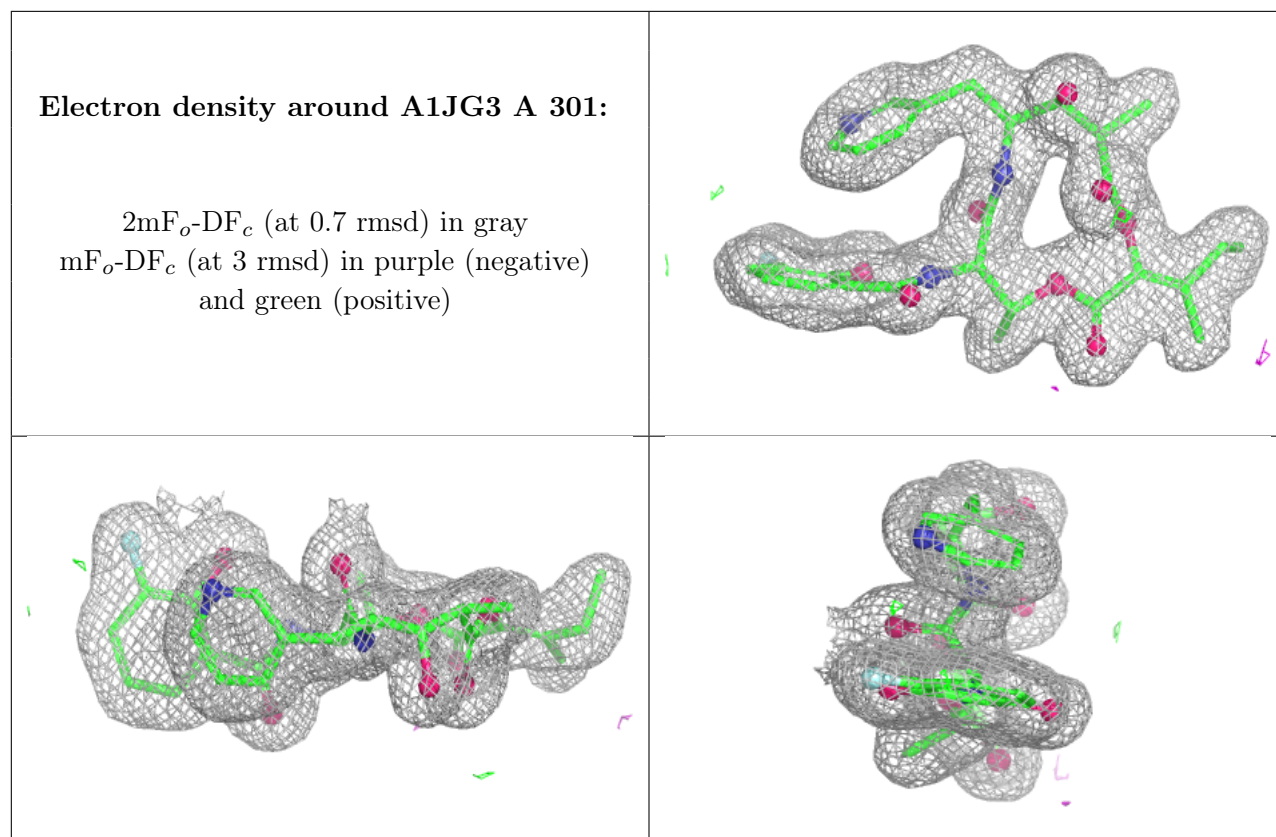
Electron density around A1JG3 C 301:

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

**Electron density around A1JG3 B 301:**

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





6.5 Other polymers [i](#)

There are no such residues in this entry.