wwPDB X-ray Structure Validation Summary Report (i)

Oct 15, 2026 – 05:22 pm BST

PDB ID	:	9T5Y / pdb_00009t5y
Title	:	Crystal structure of Mycobacterium abscessus Phosphopantetheine adenylyltransferase in complex with inhibitor
Authors	:	Thomas, S.E.; McCarthy, W.; Coyne, A.G.; Floto, R.A.
Deposited on	:	2025-11-06
Resolution	:	1.48 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i)) were used in the production of this report:

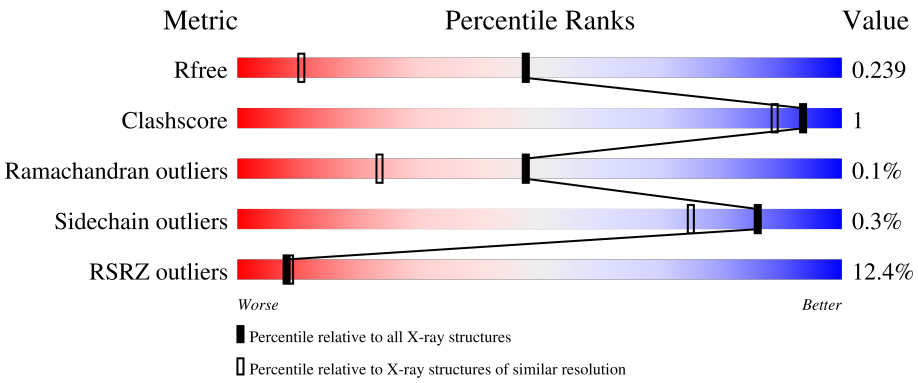
MolProbity	:	4-5-2 with Phenix2.2
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.2
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

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 1.48 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	6779 (1.50-1.46)
Clashscore	190562	7025 (1.50-1.46)
Ramachandran outliers	187476	6917 (1.50-1.46)
Sidechain outliers	187428	6914 (1.50-1.46)
RSRZ outliers	180081	6781 (1.50-1.46)

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	A	162	16% 91% • 6%
1	B	162	12% 93% • •
1	C	162	10% 91% • 5%
1	D	162	8% 95% • •
1	E	162	14% 94% • •

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Mol	Chain	Length	Quality of chain
1	F	162	11% 88% 7% 5%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 7866 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.

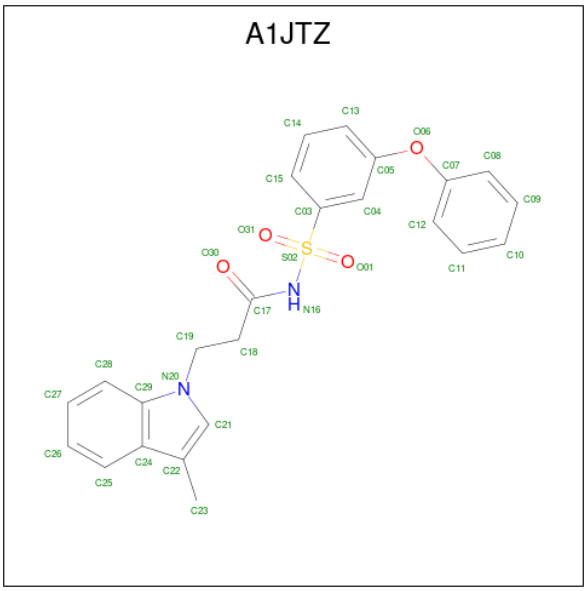
- Molecule 1 is a protein called Phosphopantetheine adenylyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	153	1166	745	197	217	7	0	5	0
1	B	156	1174	748	202	218	6	0	4	0
1	C	154	1161	737	194	224	6	0	3	0
1	D	157	1201	764	206	224	7	0	5	0
1	E	157	1186	753	201	225	7	0	3	0
1	F	154	1147	732	192	217	6	0	1	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP B1MDL6
B	0	SER	-	expression tag	UNP B1MDL6
C	0	SER	-	expression tag	UNP B1MDL6
D	0	SER	-	expression tag	UNP B1MDL6
E	0	SER	-	expression tag	UNP B1MDL6
F	0	SER	-	expression tag	UNP B1MDL6

- Molecule 2 is 3-(3-methylindol-1-yl)- {N}-(3-phenoxyphenyl)sulfonyl-propanamide (CCD ID: A1JTZ) (formula: $C_{24}H_{22}N_2O_4S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			31	24	2	4	1		
2	B	1	Total	C	N	O	S	0	0
			31	24	2	4	1		
2	C	1	Total	C	N	O	S	0	0
			31	24	2	4	1		
2	D	1	Total	C	N	O	S	0	0
			31	24	2	4	1		
2	E	1	Total	C	N	O	S	0	0
			31	24	2	4	1		
2	F	1	Total	C	N	O	S	0	0
			31	24	2	4	1		

- Molecule 3 is water.

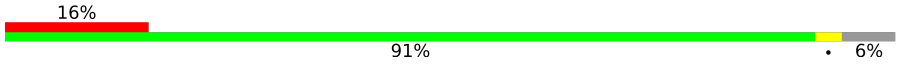
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	91	Total	O	0	0
			91	91		
3	B	125	Total	O	0	0
			125	125		
3	C	115	Total	O	0	0
			115	115		
3	D	117	Total	O	0	0
			117	117		
3	E	101	Total	O	0	0
			101	101		
3	F	96	Total	O	0	0
			96	96		

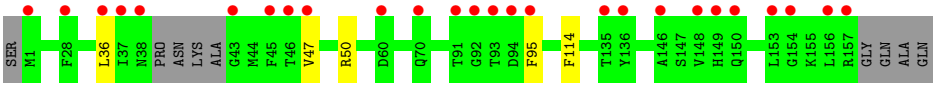
3 Residue-property plots (i)

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: Phosphopantetheine adenylyltransferase

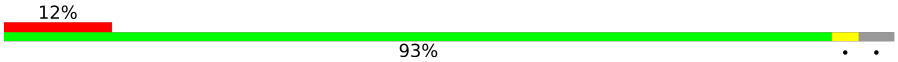
Chain A:

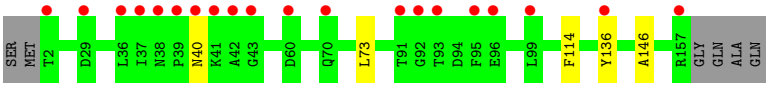




• Molecule 1: Phosphopantetheine adenylyltransferase

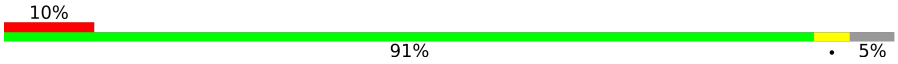
Chain B:

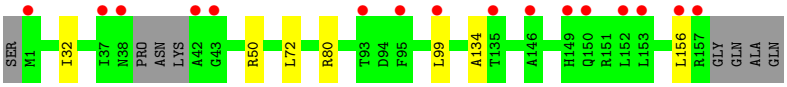




• Molecule 1: Phosphopantetheine adenylyltransferase

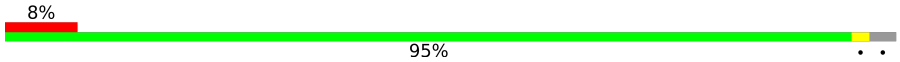
Chain C:

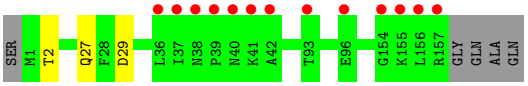




• Molecule 1: Phosphopantetheine adenylyltransferase

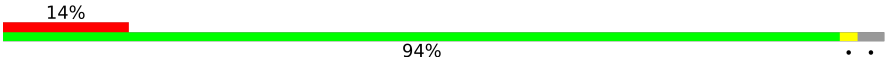
Chain D:

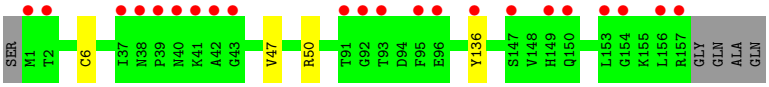




• Molecule 1: Phosphopantetheine adenylyltransferase

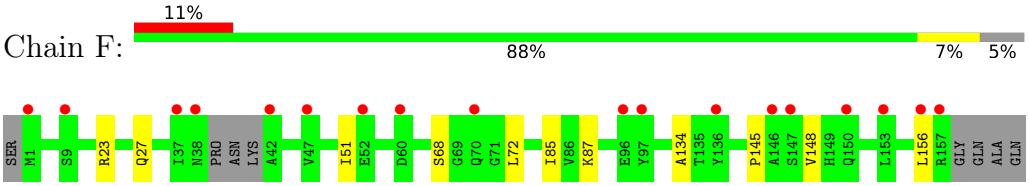
Chain E:





• Molecule 1: Phosphopantetheine adenylyltransferase

Chain F:


11%
88%
7%
5%
SER
M1
S9
R23
Q27
I37
N38
PRO
ASN
LYS
A42
V47
I51
E52
D60
S68
G69
Q70
G71
L72
I85
V86
K87
E96
Y97
A134
T135
Y136
P145
A146
S147
V148
H149
Q150
L153
L156
R157
GLY
GLN
ALA
GLN

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	67.63Å 126.38Å 132.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	63.19 – 1.48 63.19 – 1.48	Depositor EDS
% Data completeness (in resolution range)	98.7 (63.19-1.48) 98.8 (63.19-1.48)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.43 (at 1.48Å)	Xtriage
Refinement program	PHENIX (1.21.2_5419: ???)	Depositor
R, R_{free}	0.214 , 0.238 0.216 , 0.239	Depositor DCC
R_{free} test set	9428 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å^2)	22.3	Xtriage
Anisotropy	0.373	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 24.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.003 for -h,l,k	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	7866	wwPDB-VP
Average B, all atoms (Å^2)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.03% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]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.

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: A1JTZ

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	0.32	0/1199	0.57	2/1624 (0.1%)
1	B	0.28	0/1205	0.49	0/1635
1	C	0.27	0/1187	0.47	0/1607
1	D	0.27	0/1236	0.46	0/1675
1	E	0.27	0/1214	0.48	0/1646
1	F	0.26	0/1167	0.49	0/1581
All	All	0.28	0/7208	0.50	2/9768 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	95	PHE	N-CA-CB	6.35	121.21	110.49
1	A	95	PHE	CA-CB-CG	-6.34	107.46	113.80

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	1166	0	1178	4	0
1	B	1174	0	1165	4	0
1	C	1161	0	1157	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1201	0	1214	3	0
1	E	1186	0	1175	2	0
1	F	1147	0	1141	7	0
2	A	31	0	0	1	0
2	B	31	0	0	1	0
2	C	31	0	0	0	0
2	D	31	0	0	0	0
2	E	31	0	0	0	0
2	F	31	0	0	0	0
3	A	91	0	0	1	0
3	B	125	0	0	1	0
3	C	115	0	0	0	0
3	D	117	0	0	0	0
3	E	101	0	0	0	0
3	F	96	0	0	0	0
All	All	7866	0	7030	20	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.

The worst 5 of 20 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:136:TYR:HB2	1:C:72:LEU:HG	1.82	0.62
1:A:47:VAL:HG22	1:A:50[B]:ARG:HH21	1.67	0.59
3:A:301:HOH:O	1:D:27:GLN:HG2	2.05	0.56
1:B:73:LEU:HD22	2:B:201:A1JTZ:C14	2.37	0.54
1:C:32:ILE:HD13	1:C:80:ARG:HD3	1.88	0.54

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	154/162 (95%)	150 (97%)	4 (3%)	0	100	100
1	B	158/162 (98%)	153 (97%)	4 (2%)	1 (1%)	21	6
1	C	153/162 (94%)	150 (98%)	3 (2%)	0	100	100
1	D	160/162 (99%)	156 (98%)	4 (2%)	0	100	100
1	E	158/162 (98%)	154 (98%)	4 (2%)	0	100	100
1	F	151/162 (93%)	147 (97%)	4 (3%)	0	100	100
All	All	934/972 (96%)	910 (97%)	23 (2%)	1 (0%)	48	23

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	40	ASN

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	124/130 (95%)	124 (100%)	0	100	100
1	B	121/130 (93%)	121 (100%)	0	100	100
1	C	123/130 (95%)	122 (99%)	1 (1%)	73	52
1	D	128/130 (98%)	128 (100%)	0	100	100
1	E	124/130 (95%)	122 (98%)	2 (2%)	55	26
1	F	118/130 (91%)	118 (100%)	0	100	100
All	All	738/780 (95%)	735 (100%)	3 (0%)	86	70

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

Mol	Chain	Res	Type
1	C	99	LEU
1	E	6[A]	CYS
1	E	6[B]	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	70	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 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	A1JTZ	F	201	-	34,34,34	1.17	3 (8%)	46,48,48	2.69	8 (17%)
2	A1JTZ	E	201	-	34,34,34	1.10	4 (11%)	46,48,48	2.66	10 (21%)
2	A1JTZ	C	201	-	34,34,34	1.10	3 (8%)	46,48,48	2.54	8 (17%)
2	A1JTZ	B	201	-	34,34,34	1.15	3 (8%)	46,48,48	2.72	9 (19%)
2	A1JTZ	D	201	-	34,34,34	1.12	3 (8%)	46,48,48	2.73	8 (17%)
2	A1JTZ	A	201	-	34,34,34	1.15	4 (11%)	46,48,48	2.69	8 (17%)

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	A1JTZ	F	201	-	-	1/20/20/20	0/4/4/4
2	A1JTZ	E	201	-	-	0/20/20/20	0/4/4/4
2	A1JTZ	C	201	-	-	2/20/20/20	0/4/4/4
2	A1JTZ	B	201	-	-	3/20/20/20	0/4/4/4
2	A1JTZ	D	201	-	-	0/20/20/20	0/4/4/4
2	A1JTZ	A	201	-	-	0/20/20/20	0/4/4/4

The worst 5 of 20 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	201	A1JTZ	C21-C22	3.16	1.40	1.35
2	F	201	A1JTZ	C21-C22	3.04	1.39	1.35
2	D	201	A1JTZ	C21-C22	3.01	1.39	1.35
2	A	201	A1JTZ	C21-C22	2.97	1.39	1.35
2	B	201	A1JTZ	C21-C22	2.96	1.39	1.35

The worst 5 of 51 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	F	201	A1JTZ	O31-S02-O01	-14.19	102.10	119.55
2	A	201	A1JTZ	O31-S02-O01	-14.06	102.27	119.55
2	B	201	A1JTZ	O31-S02-O01	-13.95	102.40	119.55
2	E	201	A1JTZ	O31-S02-O01	-13.61	102.82	119.55
2	C	201	A1JTZ	O31-S02-O01	-13.36	103.12	119.55

There are no chirality outliers.

5 of 6 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	201	A1JTZ	C17-N16-S02-C03
2	B	201	A1JTZ	C17-N16-S02-O31
2	B	201	A1JTZ	C17-N16-S02-O01
2	C	201	A1JTZ	N16-C17-C18-C19
2	C	201	A1JTZ	O30-C17-C18-C19

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	201	A1JTZ	1	0

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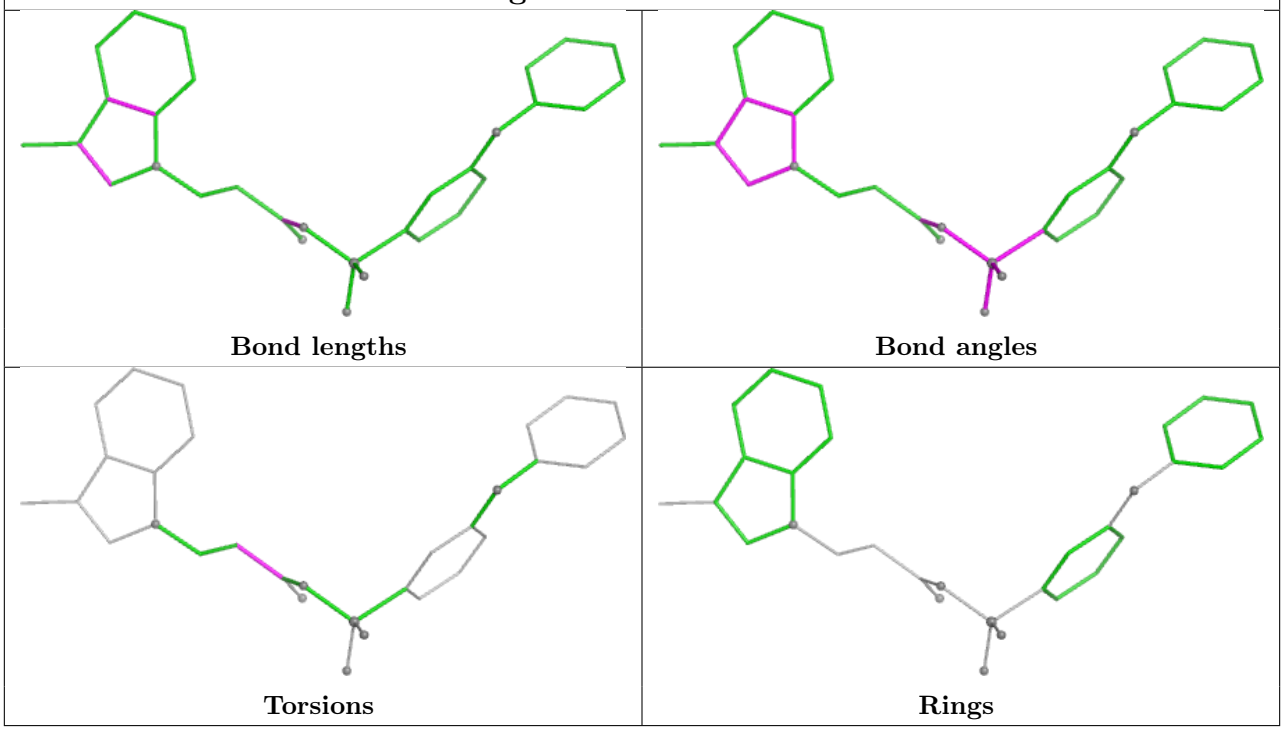
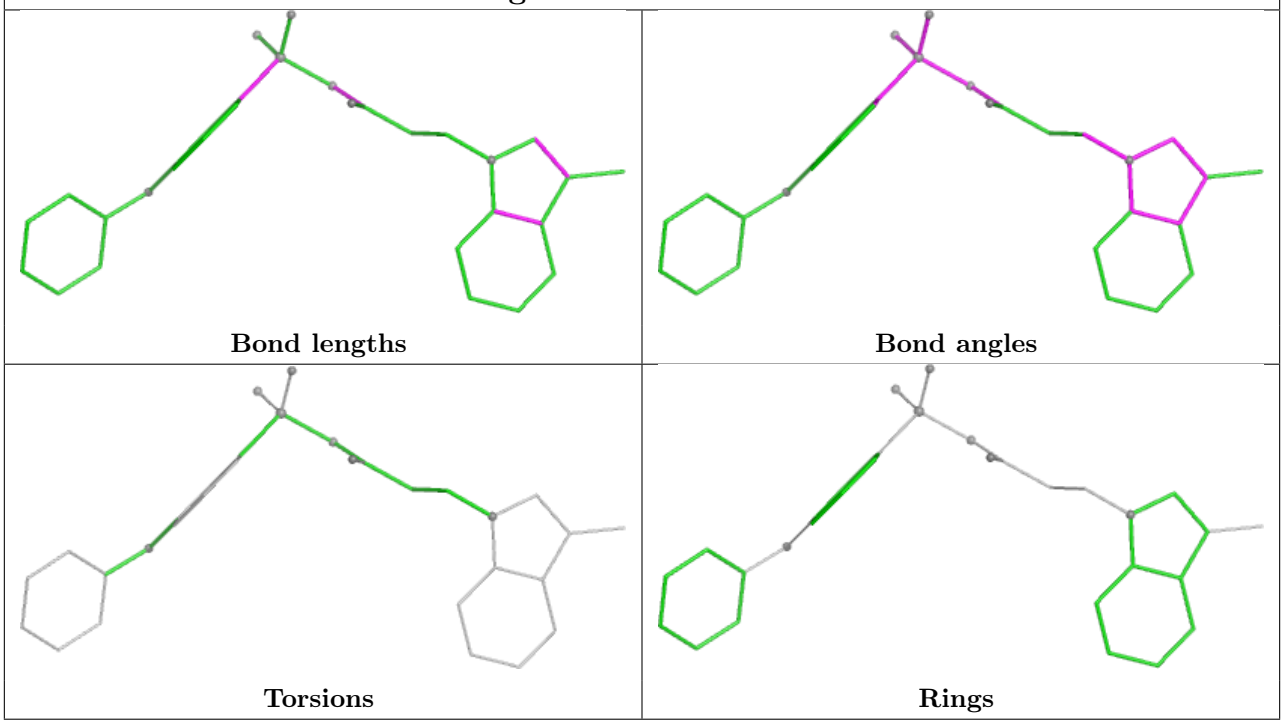
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	201	A1JTZ	1	0

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.

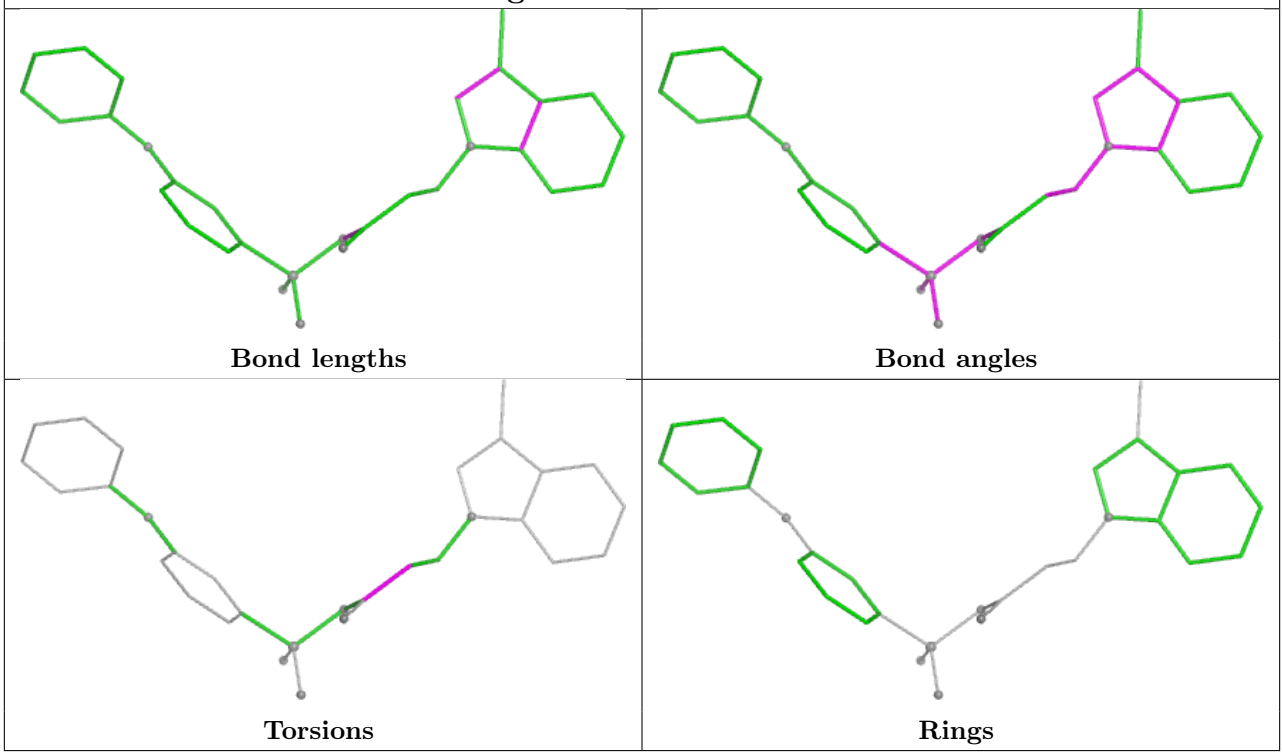
Ligand A1JTZ F 201



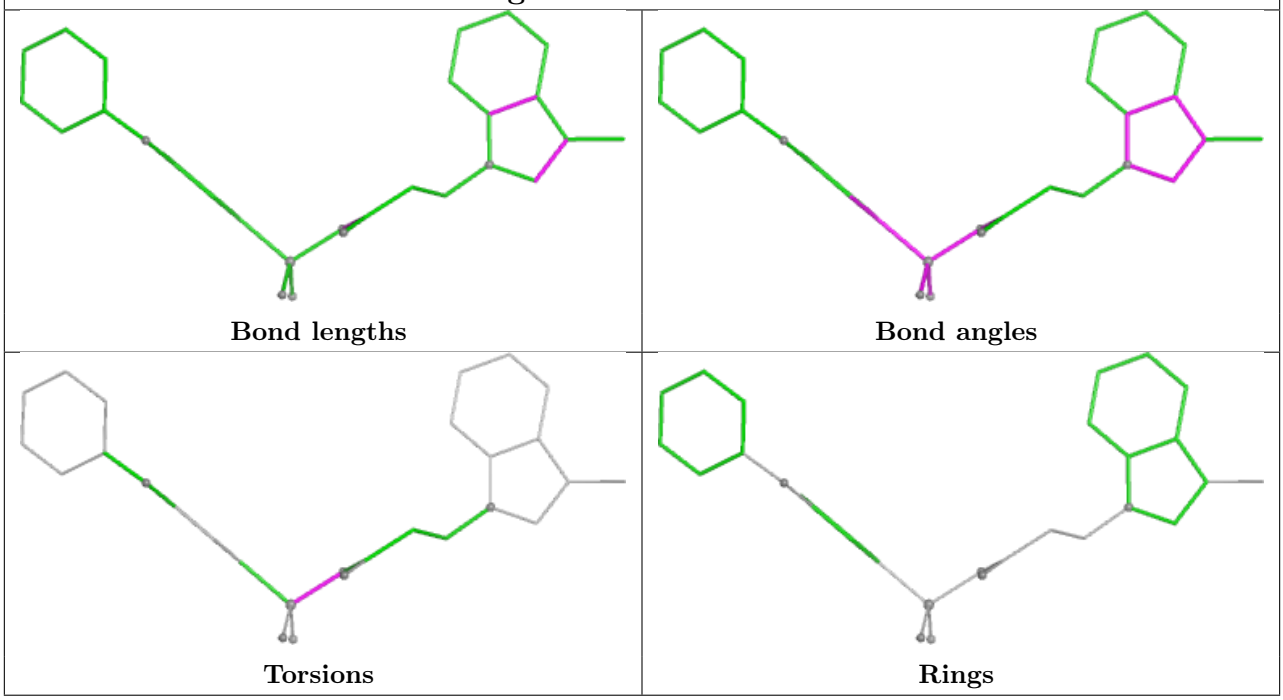
Ligand A1JTZ E 201



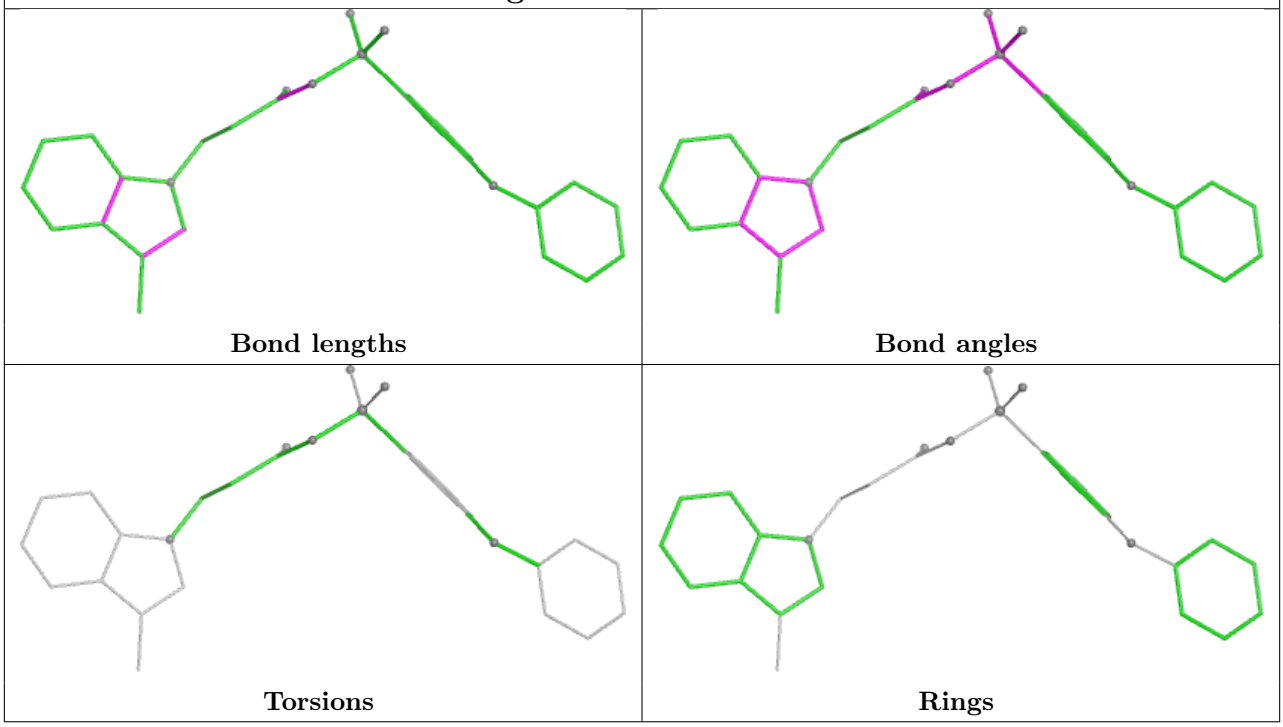
Ligand A1JTZ C 201

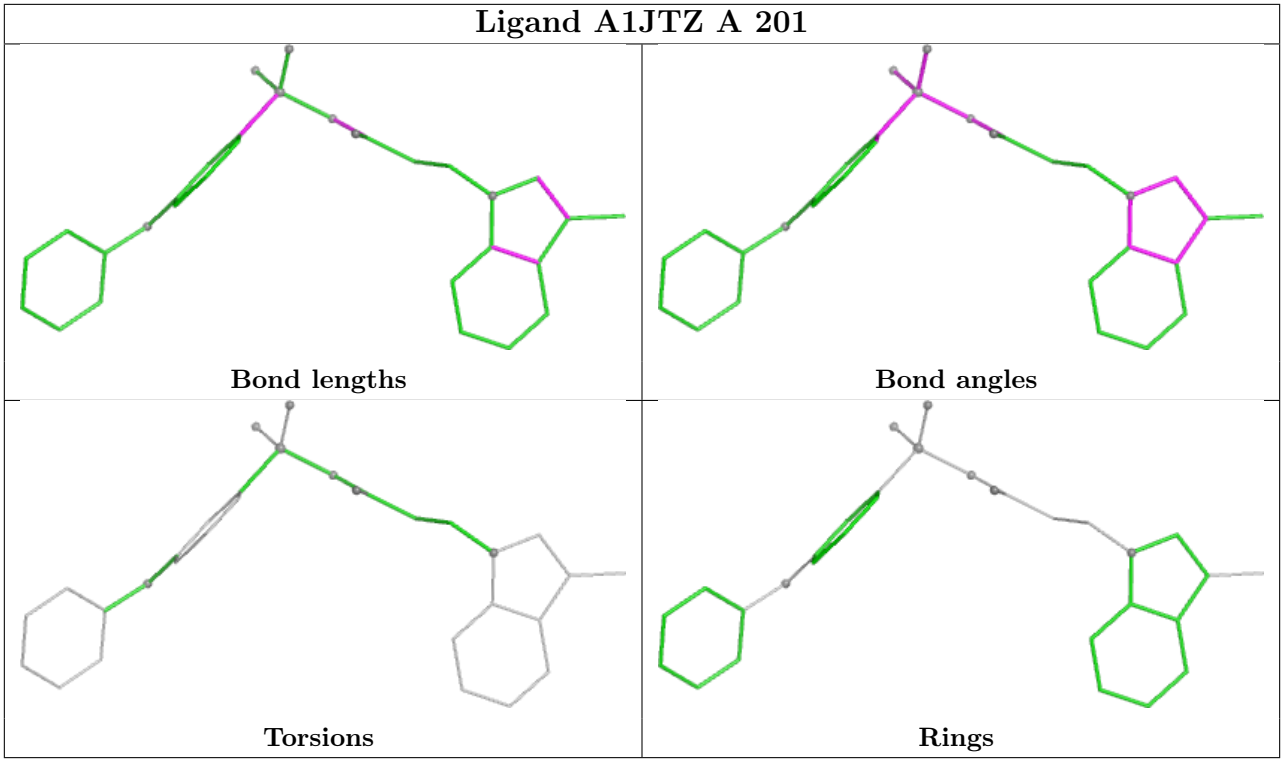


Ligand A1JTZ B 201



Ligand A1JTZ D 201





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, 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	153/162 (94%)	0.96	26 (16%)	4	4	13, 27, 49, 80	5 (3%)
1	B	156/162 (96%)	0.78	20 (12%)	7	8	12, 24, 45, 75	4 (2%)
1	C	154/162 (95%)	0.66	16 (10%)	11	12	16, 25, 43, 60	3 (1%)
1	D	157/162 (96%)	0.74	13 (8%)	17	19	14, 24, 49, 114	5 (3%)
1	E	157/162 (96%)	0.72	22 (14%)	6	7	12, 24, 46, 75	3 (1%)
1	F	154/162 (95%)	0.77	18 (11%)	9	10	18, 26, 47, 62	1 (0%)
All	All	931/972 (95%)	0.77	115 (12%)	8	9	12, 25, 47, 114	21 (2%)

The worst 5 of 115 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	156	LEU	8.0
1	A	38	ASN	7.6
1	B	42	ALA	7.5
1	C	42	ALA	7.3
1	B	91	THR	7.2

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 oligosaccharides in this entry.

6.4 Ligands (i)

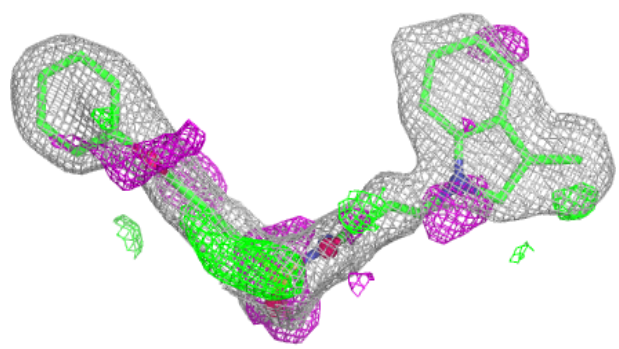
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

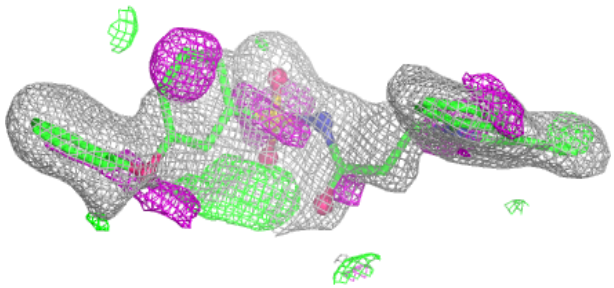
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	A1JTZ	B	201	31/31	0.72	0.17	29,39,58,71	0
2	A1JTZ	A	201	31/31	0.73	0.17	31,41,61,67	0
2	A1JTZ	C	201	31/31	0.76	0.17	30,37,49,59	0
2	A1JTZ	E	201	31/31	0.78	0.15	27,36,52,59	0
2	A1JTZ	D	201	31/31	0.80	0.15	29,39,58,67	0
2	A1JTZ	F	201	31/31	0.82	0.15	28,35,53,60	0

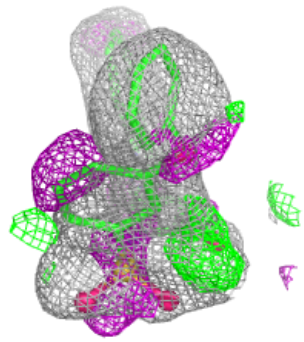
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around A1JTZ B 201:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative) and green (positive)

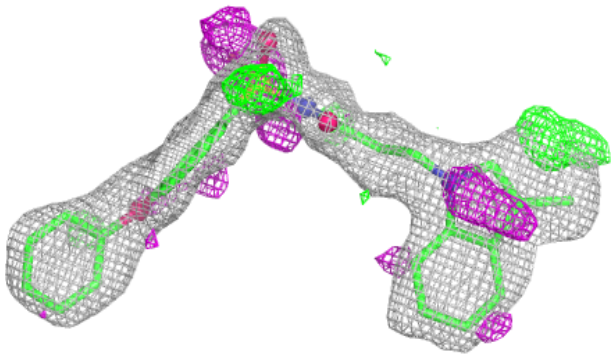


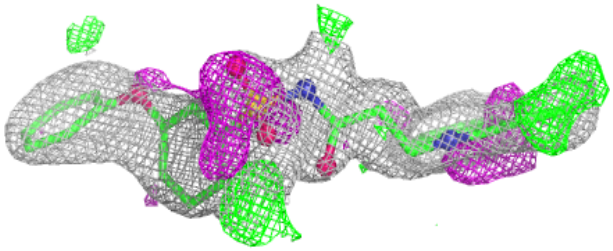


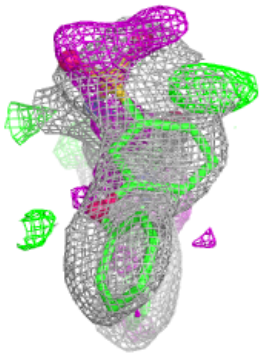


Electron density around A1JTZ A 201:

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

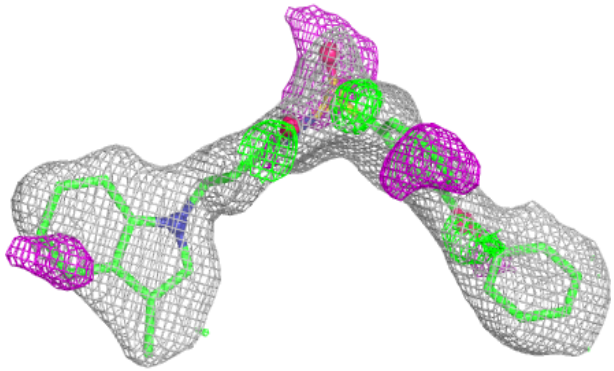


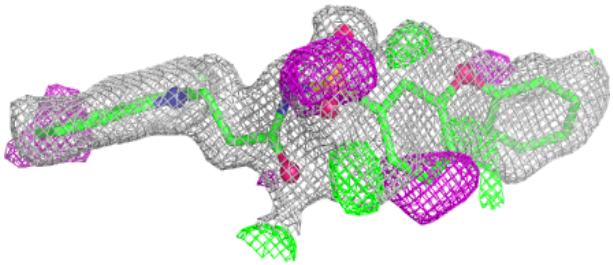


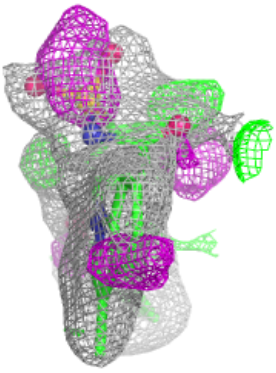


Electron density around A1JTZ C 201:

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

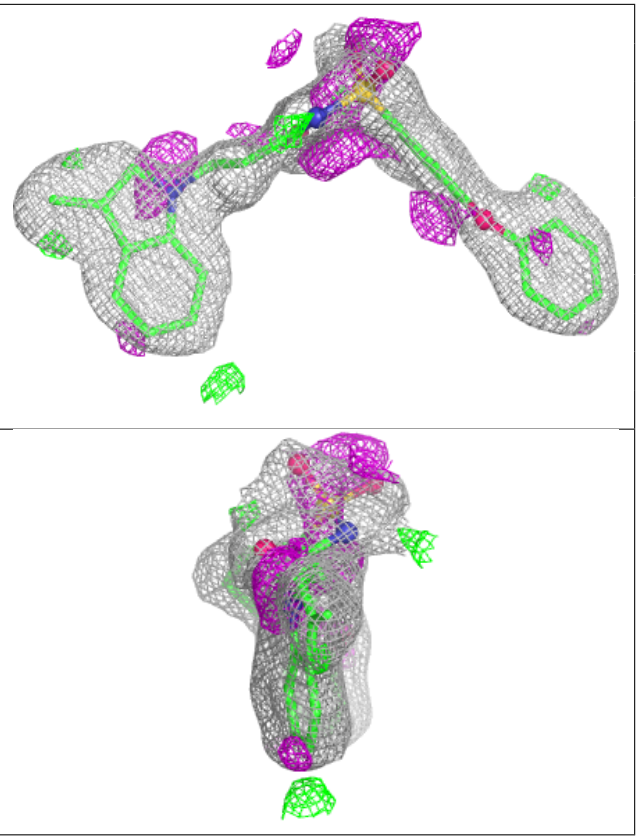






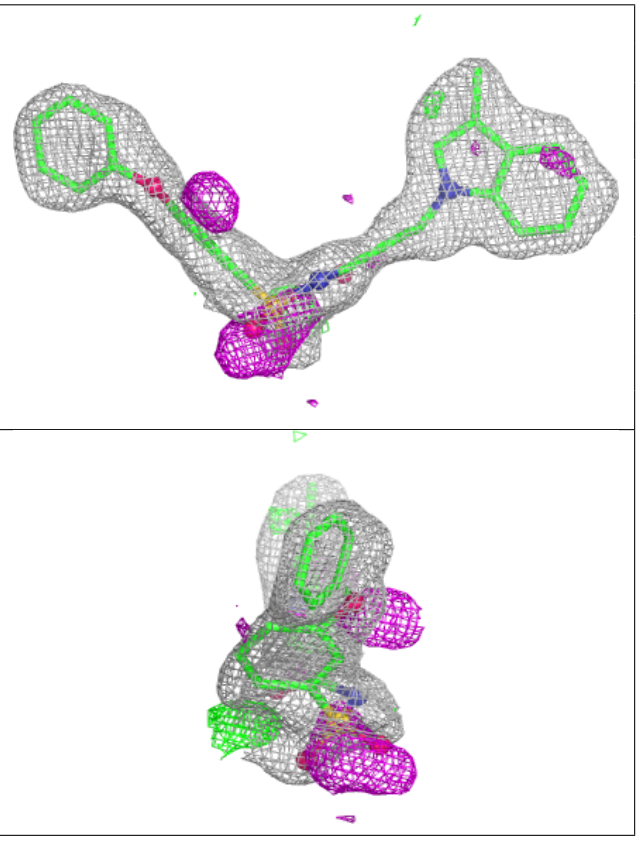
Electron density around A1JTZ E 201:

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



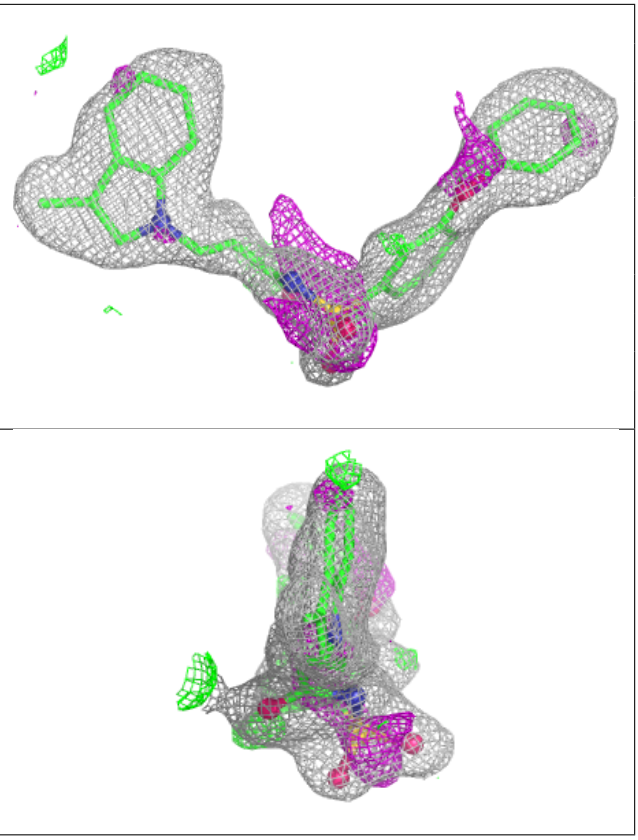
Electron density around A1JTZ D 201:

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



Electron density around A1JTZ F 201:

$2mF_o$-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.