Full wwPDB X-ray Structure Validation Report

Aug 30, 2022 – 04:51 pm BST

PDB ID	:	7PKK
Title	:	LpxC Inhibitors With Fluoroproline As A Novel Zinc-Binding Group Can Serve As A Novel Class of Antibiotic With Activity Against Multidrug-Resistant Gram-Negative Bacteria
Authors	:	Ryan, M.D.; Pallin, T.D.; Lamers, M.B.A.C.; Leonard, P.M.
Deposited on	:	2021-08-25
Resolution	:	2.55 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:

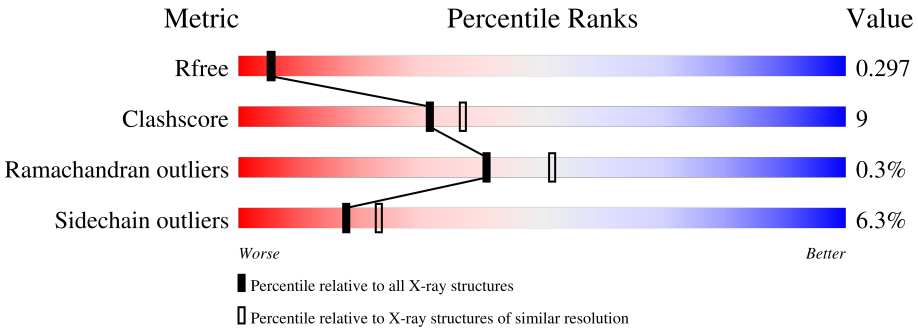
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.30
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0267
CCP4	:	7.1.010 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.30

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 2.55 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	1284 (2.56-2.52)
Clashscore	141614	1332 (2.56-2.52)
Ramachandran outliers	138981	1315 (2.56-2.52)
Sidechain outliers	138945	1315 (2.56-2.52)

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%

Mol	Chain	Length	Quality of chain
1	A	312	74% 20% • 5%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 2410 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 UDP-3-O-acyl-N-acetylglucosamine deacetylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	297	Total	C	N	O	S	0	1	0
			2264	1441	380	439	4			

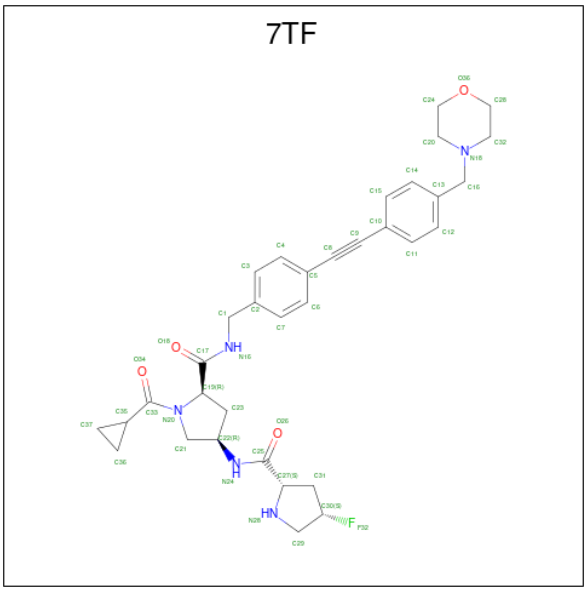
There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-12	HIS	-	expression tag	UNP P47205
A	-11	HIS	-	expression tag	UNP P47205
A	-10	HIS	-	expression tag	UNP P47205
A	-9	HIS	-	expression tag	UNP P47205
A	-8	HIS	-	expression tag	UNP P47205
A	-7	HIS	-	expression tag	UNP P47205
A	-6	GLU	-	expression tag	UNP P47205
A	-5	ASN	-	expression tag	UNP P47205
A	-4	LEU	-	expression tag	UNP P47205
A	-3	TYR	-	expression tag	UNP P47205
A	-2	PHE	-	expression tag	UNP P47205
A	-1	GLN	-	expression tag	UNP P47205
A	0	SER	-	expression tag	UNP P47205
A	40	SER	CYS	conflict	UNP P47205

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is (2R,4R)-1-cyclopropylcarbonyl-4-[[(2S,4S)-4-fluoranylpyrrolidin-2-yl]carbonyl amino]-N-[[4-[2-[4-(morpholin-4-ylmethyl)phenyl]ethynyl]phenyl]methyl]pyrrolidine-2-carb oxamide (three-letter code: 7TF) (formula: $C_{34}H_{40}FN_5O_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			44	34	1	5	4		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	101	Total	O	0	0
			101	101		

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. 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. 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: UDP-3-O-acyl-N-acetylglucosamine deacetylase

Chain A:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	35.55Å 66.96Å 62.96Å 90.00° 90.55° 90.00°	Depositor
Resolution (Å)	35.55 – 2.55 35.55 – 2.55	Depositor EDS
% Data completeness (in resolution range)	98.0 (35.55-2.55) 98.0 (35.55-2.55)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.34 (at 2.54Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.202 , 0.297 0.210 , 0.297	Depositor DCC
R_{free} test set	478 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å^2)	51.9	Xtriage
Anisotropy	0.143	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	(Not available) , (Not available)	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.049 for h,-k,-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	2410	wwPDB-VP
Average B, all atoms (Å^2)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.96% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, 7TF

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	0.95	1/2305 (0.0%)	0.94	0/3122

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	40	SER	CB-OG	29.75	1.80	1.42

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	2264	0	2228	41	0
2	A	1	0	0	0	0
3	A	44	0	0	0	0
4	A	101	0	0	9	0
All	All	2410	0	2228	41	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 9.

All (41) 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:40:SER:OG	1:A:40:SER:CB	1.80	1.29
1:A:272:ARG:NH2	4:A:601:HOH:O	1.88	0.99
1:A:99:GLU:OE1	4:A:602:HOH:O	2.04	0.75
1:A:227:GLY:O	4:A:603:HOH:O	2.12	0.66
1:A:174:VAL:HG13	1:A:179:THR:HG21	1.78	0.66
1:A:88:ILE:HD11	1:A:114:ILE:HG21	1.82	0.61
1:A:193:PHE:O	1:A:197:ILE:HD12	2.02	0.60
1:A:169:THR:O	4:A:604:HOH:O	2.17	0.58
1:A:11:ILE:HG21	1:A:113:LEU:HB3	1.86	0.58
1:A:256:GLU:OE2	1:A:258:ARG:NE	2.27	0.54
1:A:165:PHE:HB3	1:A:170:GLN:NE2	2.25	0.51
1:A:269:GLN:HG2	4:A:678:HOH:O	2.10	0.51
1:A:175:ASP:OD1	1:A:175:ASP:C	2.50	0.50
1:A:146:PHE:CE2	1:A:274:LEU:HD22	2.46	0.50
1:A:268:ASN:O	1:A:272:ARG:HG3	2.11	0.50
1:A:130:ARG:HG3	1:A:254:ILE:HB	1.94	0.50
1:A:104:ASP:C	1:A:213:ASN:HA	2.34	0.48
1:A:154:VAL:HG11	1:A:239:ILE:HG23	1.96	0.48
1:A:99:GLU:OE2	4:A:605:HOH:O	2.20	0.47
1:A:162:HIS:HB3	1:A:165:PHE:HD1	1.80	0.47
1:A:165:PHE:HB3	1:A:170:GLN:CD	2.35	0.46
1:A:108:GLY:N	1:A:109:PRO:CD	2.79	0.45
1:A:174:VAL:HG13	1:A:179:THR:CG2	2.43	0.45
1:A:8:LYS:HE2	4:A:651:HOH:O	2.16	0.45
1:A:152:PHE:CE1	1:A:253:LEU:HD23	2.52	0.45
1:A:62:MET:CB	4:A:674:HOH:O	2.65	0.44
1:A:64:THR:HG1	1:A:264:HIS:CE1	2.36	0.44
1:A:73:VAL:HA	1:A:98:SER:O	2.19	0.43
1:A:146:PHE:CD2	1:A:274:LEU:HD22	2.53	0.43
1:A:156:PHE:HA	1:A:259:GLY:O	2.19	0.43
1:A:152:PHE:HE1	1:A:253:LEU:HD23	1.84	0.43
1:A:100:VAL:HG13	1:A:101:PRO:HD2	2.02	0.42
1:A:131:ILE:HD11	1:A:253:LEU:HD11	2.01	0.42
1:A:179:THR:HG23	1:A:184:GLU:HG3	2.02	0.41
1:A:132:LYS:HB2	1:A:280:ALA:O	2.20	0.41
1:A:19:HIS:HB2	4:A:602:HOH:O	2.20	0.41
1:A:191:PHE:HB3	1:A:215:ILE:HB	2.02	0.41
1:A:64:THR:OG1	1:A:264:HIS:ND1	2.49	0.41
1:A:76:VAL:HG23	1:A:80:LEU:HD12	2.02	0.41
1:A:122:GLN:O	1:A:124:ALA:N	2.54	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	294/312 (94%)	273 (93%)	20 (7%)	1 (0%)	41 51

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	123	GLU

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	240/266 (90%)	225 (94%)	15 (6%)	18 23

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	ILE
1	A	55	GLU
1	A	116	SER
1	A	134	GLU
1	A	136	SER
1	A	150	ASP
1	A	153	LYS

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Mol	Chain	Res	Type
1	A	156	PHE
1	A	169	THR
1	A	171	GLN
1	A	178	SER
1	A	197	ILE
1	A	221	ARG
1	A	246	LEU
1	A	247	TYR

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

Mol	Chain	Res	Type
1	A	171	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
3	7TF	A	502	2	49,49,49	1.30	2 (4%)	61,68,68	0.92	2 (3%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	7TF	A	502	2	-	2/34/65/65	0/6/6/6

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502	7TF	C10-C9	-5.93	1.30	1.44
3	A	502	7TF	C5-C8	-5.81	1.30	1.44

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	502	7TF	C37-C35-C33	-2.68	113.53	117.01
3	A	502	7TF	C16-N18-C32	-2.45	105.61	111.06

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	7TF	O26-C25-C27-C31
3	A	502	7TF	N24-C25-C27-C31

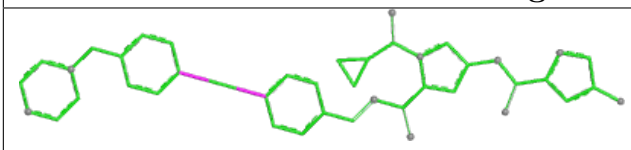
There are no ring outliers.

No monomer is involved in short contacts.

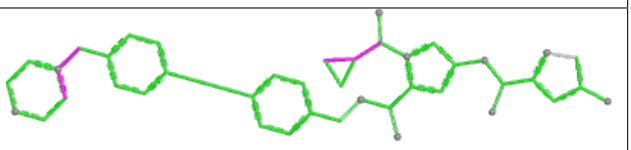
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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.

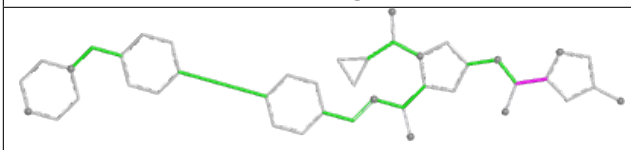
Ligand 7TF A 502



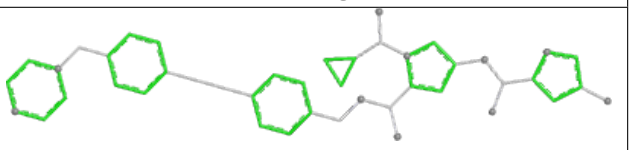
Bond lengths



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

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

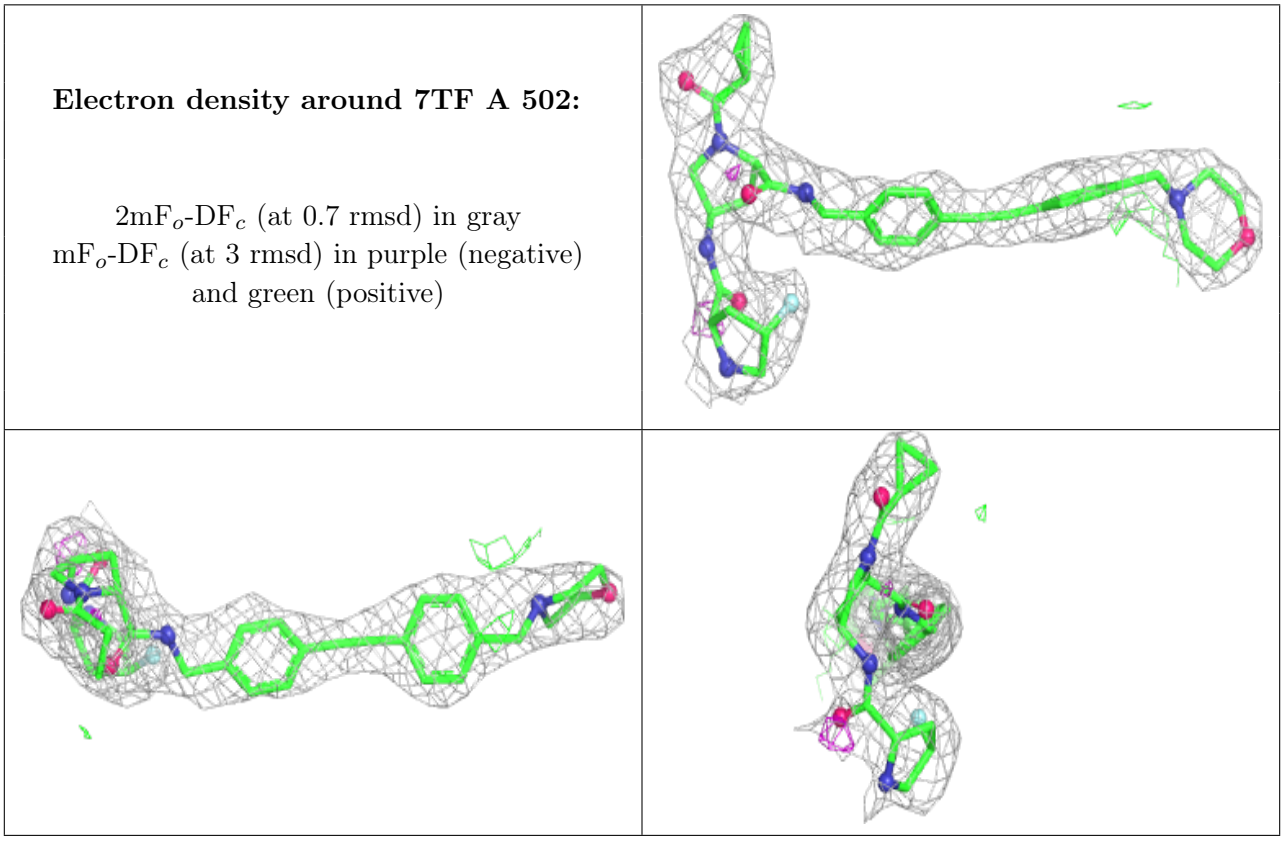
6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

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 7TF A 502:

$2mF_o$-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)

Unable to reproduce the depositors R factor - this section is therefore empty.