



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 28, 2026 – 08:53 PM UTC

PDB ID : 6VKO / pdb_00006vko
Title : Crystal Structure of human PARP-1 CAT domain bound to inhibitor UKTT15
Authors : Langelier, M.F.; Pascal, J.M.
Deposited on : 2020-01-21
Resolution : 2.80 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 ⓘ 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

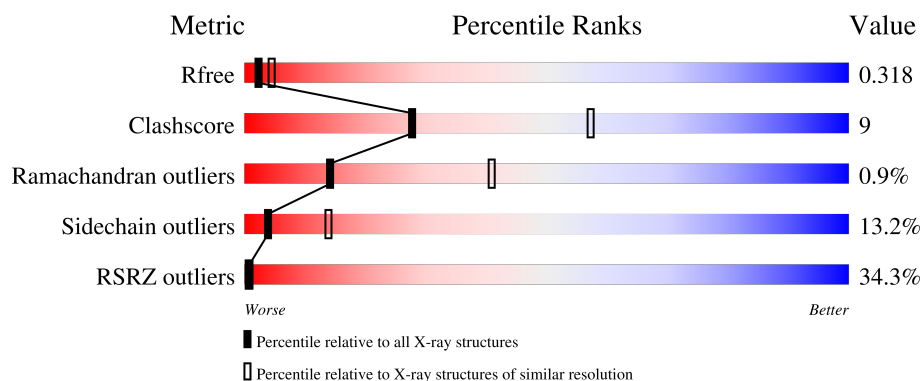
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	372	
1	B	372	
1	C	372	
1	D	372	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10970 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 Poly [ADP-ribose] polymerase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	340	Total	C	N	O	S	0	1	0
			2686	1713	455	507	11			
1	B	340	Total	C	N	O	S	0	1	0
			2683	1712	454	506	11			
1	C	340	Total	C	N	O	S	0	1	0
			2683	1712	454	506	11			
1	D	340	Total	C	N	O	S	0	1	0
			2683	1712	454	506	11			

There are 88 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	640	MET	-	initiating methionine	UNP P09874
A	641	GLY	-	expression tag	UNP P09874
A	642	SER	-	expression tag	UNP P09874
A	643	SER	-	expression tag	UNP P09874
A	644	HIS	-	expression tag	UNP P09874
A	645	HIS	-	expression tag	UNP P09874
A	646	HIS	-	expression tag	UNP P09874
A	647	HIS	-	expression tag	UNP P09874
A	648	HIS	-	expression tag	UNP P09874
A	649	HIS	-	expression tag	UNP P09874
A	650	SER	-	expression tag	UNP P09874
A	651	SER	-	expression tag	UNP P09874
A	652	GLY	-	expression tag	UNP P09874
A	653	LEU	-	expression tag	UNP P09874
A	654	VAL	-	expression tag	UNP P09874
A	655	PRO	-	expression tag	UNP P09874
A	656	ARG	-	expression tag	UNP P09874
A	657	GLY	-	expression tag	UNP P09874
A	658	SER	-	expression tag	UNP P09874
A	659	HIS	-	expression tag	UNP P09874
A	660	MET	-	expression tag	UNP P09874

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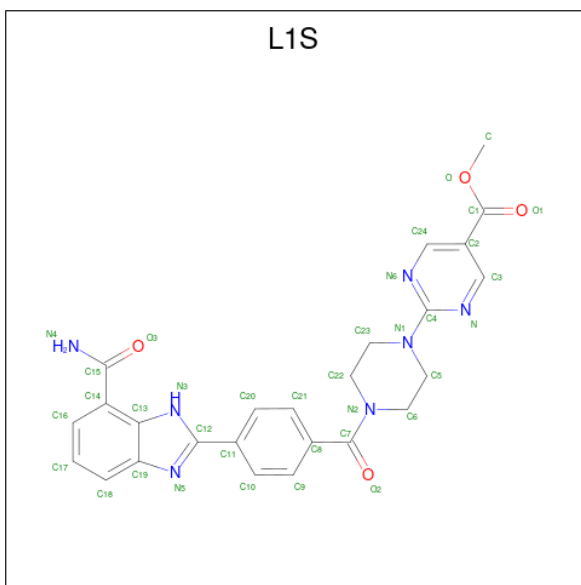
Chain	Residue	Modelled	Actual	Comment	Reference
A	762	ALA	VAL	variant	UNP P09874
B	640	MET	-	initiating methionine	UNP P09874
B	641	GLY	-	expression tag	UNP P09874
B	642	SER	-	expression tag	UNP P09874
B	643	SER	-	expression tag	UNP P09874
B	644	HIS	-	expression tag	UNP P09874
B	645	HIS	-	expression tag	UNP P09874
B	646	HIS	-	expression tag	UNP P09874
B	647	HIS	-	expression tag	UNP P09874
B	648	HIS	-	expression tag	UNP P09874
B	649	HIS	-	expression tag	UNP P09874
B	650	SER	-	expression tag	UNP P09874
B	651	SER	-	expression tag	UNP P09874
B	652	GLY	-	expression tag	UNP P09874
B	653	LEU	-	expression tag	UNP P09874
B	654	VAL	-	expression tag	UNP P09874
B	655	PRO	-	expression tag	UNP P09874
B	656	ARG	-	expression tag	UNP P09874
B	657	GLY	-	expression tag	UNP P09874
B	658	SER	-	expression tag	UNP P09874
B	659	HIS	-	expression tag	UNP P09874
B	660	MET	-	expression tag	UNP P09874
B	762	ALA	VAL	variant	UNP P09874
C	640	MET	-	initiating methionine	UNP P09874
C	641	GLY	-	expression tag	UNP P09874
C	642	SER	-	expression tag	UNP P09874
C	643	SER	-	expression tag	UNP P09874
C	644	HIS	-	expression tag	UNP P09874
C	645	HIS	-	expression tag	UNP P09874
C	646	HIS	-	expression tag	UNP P09874
C	647	HIS	-	expression tag	UNP P09874
C	648	HIS	-	expression tag	UNP P09874
C	649	HIS	-	expression tag	UNP P09874
C	650	SER	-	expression tag	UNP P09874
C	651	SER	-	expression tag	UNP P09874
C	652	GLY	-	expression tag	UNP P09874
C	653	LEU	-	expression tag	UNP P09874
C	654	VAL	-	expression tag	UNP P09874
C	655	PRO	-	expression tag	UNP P09874
C	656	ARG	-	expression tag	UNP P09874
C	657	GLY	-	expression tag	UNP P09874
C	658	SER	-	expression tag	UNP P09874

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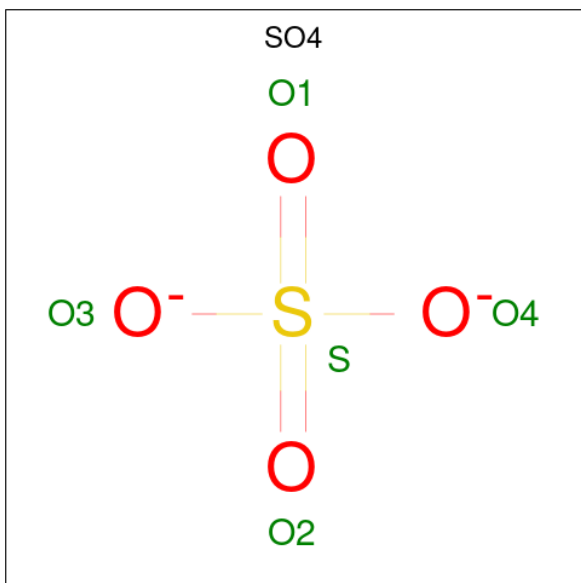
Chain	Residue	Modelled	Actual	Comment	Reference
C	659	HIS	-	expression tag	UNP P09874
C	660	MET	-	expression tag	UNP P09874
C	762	ALA	VAL	variant	UNP P09874
D	640	MET	-	initiating methionine	UNP P09874
D	641	GLY	-	expression tag	UNP P09874
D	642	SER	-	expression tag	UNP P09874
D	643	SER	-	expression tag	UNP P09874
D	644	HIS	-	expression tag	UNP P09874
D	645	HIS	-	expression tag	UNP P09874
D	646	HIS	-	expression tag	UNP P09874
D	647	HIS	-	expression tag	UNP P09874
D	648	HIS	-	expression tag	UNP P09874
D	649	HIS	-	expression tag	UNP P09874
D	650	SER	-	expression tag	UNP P09874
D	651	SER	-	expression tag	UNP P09874
D	652	GLY	-	expression tag	UNP P09874
D	653	LEU	-	expression tag	UNP P09874
D	654	VAL	-	expression tag	UNP P09874
D	655	PRO	-	expression tag	UNP P09874
D	656	ARG	-	expression tag	UNP P09874
D	657	GLY	-	expression tag	UNP P09874
D	658	SER	-	expression tag	UNP P09874
D	659	HIS	-	expression tag	UNP P09874
D	660	MET	-	expression tag	UNP P09874
D	762	ALA	VAL	variant	UNP P09874

- Molecule 2 is methyl 2-{4-[4-(7-carbamoyl-1H-benzimidazol-2-yl)benzene-1-carbonyl]piperazin-1-yl}pyrimidine-5-carboxylate (CCD ID: L1S) (formula: C₂₅H₂₃N₇O₄) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			36	25	7	4		
2	B	1	Total	C	N	O	0	0
			36	25	7	4		
2	C	1	Total	C	N	O	0	0
			36	25	7	4		
2	D	1	Total	C	N	O	0	0
			36	25	7	4		

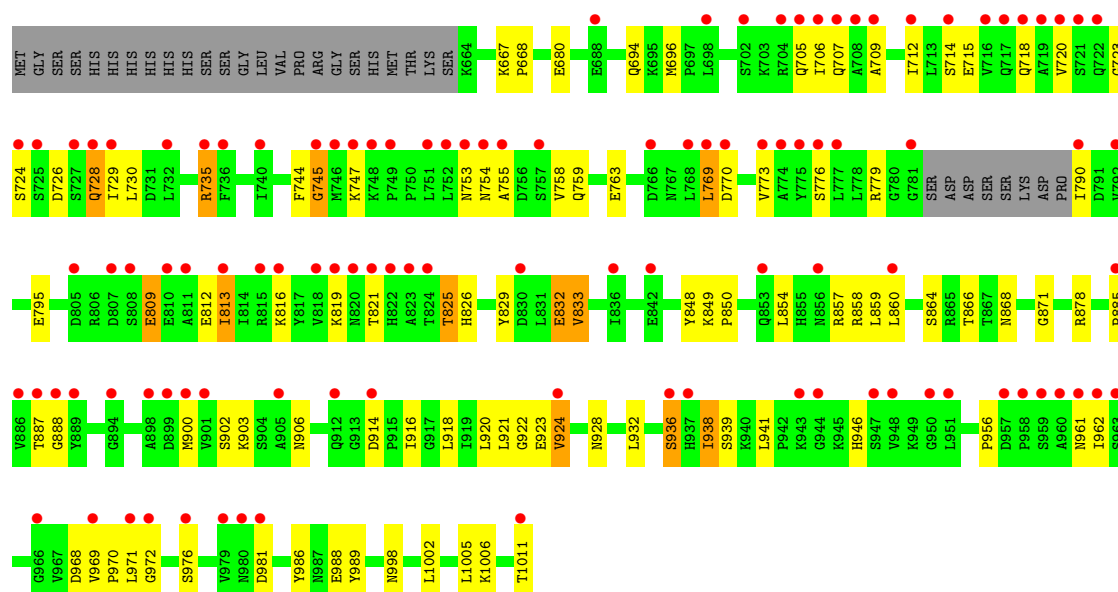
- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	6	Total O 6 6	0	0
4	B	8	Total O 8 8	0	0
4	C	7	Total O 7 7	0	0
4	D	10	Total O 10 10	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	139.70Å 129.75Å 102.72Å 90.00° 111.36° 90.00°	Depositor
Resolution (Å)	20.00 – 2.80 20.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	94.7 (20.00-2.80) 94.7 (20.00-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.16 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.8.0257	Depositor
R, R_{free}	0.284 , 0.312 0.290 , 0.318	Depositor DCC
R_{free} test set	2030 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	51.0	Xtriage
Anisotropy	0.294	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 76.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	10970	wwPDB-VP
Average B, all atoms (Å ²)	98.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 61.30 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.3371e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: LIS, SO4

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	1.06	0/2736	1.51	3/3692 (0.1%)
1	B	1.06	0/2736	1.50	4/3692 (0.1%)
1	C	1.07	1/2736 (0.0%)	1.50	3/3692 (0.1%)
1	D	1.06	0/2736	1.51	3/3692 (0.1%)
All	All	1.06	1/10944 (0.0%)	1.51	13/14768 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	C	0	2
1	D	0	1
All	All	0	5

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	963	SER	C-O	5.87	1.30	1.24

The worst 5 of 13 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	885	PRO	N-CA-C	7.90	123.68	111.11
1	B	885	PRO	N-CA-C	7.86	123.61	111.11
1	C	885	PRO	N-CA-C	7.84	123.58	111.11
1	D	885	PRO	N-CA-C	7.67	123.31	111.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	885	PRO	CA-C-N	6.21	128.82	120.50

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	920	LEU	Peptide
1	B	920	LEU	Peptide
1	C	887	THR	Peptide
1	C	920	LEU	Peptide
1	D	920	LEU	Peptide

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	2686	0	2730	48	0
1	B	2683	0	2731	43	0
1	C	2683	0	2731	53	0
1	D	2683	0	2731	53	0
2	A	36	0	0	0	0
2	B	36	0	0	0	0
2	C	36	0	0	3	0
2	D	36	0	0	4	0
3	A	15	0	0	1	0
3	B	15	0	0	2	0
3	C	15	0	0	2	0
3	D	15	0	0	2	0
4	A	6	0	0	0	1
4	B	8	0	0	0	0
4	C	7	0	0	0	0
4	D	10	0	0	0	0
All	All	10970	0	10923	197	1

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.

The worst 5 of 197 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:858:ARG:NH2	3:B:1103:SO4:O4	1.90	1.05
1:C:858:ARG:NH2	3:C:1103:SO4:O3	1.93	1.01
1:B:813:ILE:HD12	1:B:813:ILE:O	1.83	0.79
1:C:813:ILE:HD12	1:C:813:ILE:O	1.83	0.78
1:A:813:ILE:O	1:A:813:ILE:HD12	1.83	0.77

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:1204:HOH:O	4:A:1204:HOH:O[2_657]	1.35	0.85

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	337/372 (91%)	307 (91%)	27 (8%)	3 (1%)	14	41
1	B	337/372 (91%)	309 (92%)	25 (7%)	3 (1%)	14	41
1	C	337/372 (91%)	306 (91%)	28 (8%)	3 (1%)	14	41
1	D	337/372 (91%)	307 (91%)	27 (8%)	3 (1%)	14	41
All	All	1348/1488 (91%)	1229 (91%)	107 (8%)	12 (1%)	14	41

5 of 12 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	723	GLY
1	A	745	GLY
1	B	723	GLY
1	B	745	GLY
1	C	723	GLY

5.3.2 Protein sidechains ⓘ

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	298/326 (91%)	258 (87%)	40 (13%)	4	13
1	B	298/326 (91%)	258 (87%)	40 (13%)	4	13
1	C	298/326 (91%)	260 (87%)	38 (13%)	4	15
1	D	298/326 (91%)	259 (87%)	39 (13%)	4	14
All	All	1192/1304 (91%)	1035 (87%)	157 (13%)	4	14

5 of 157 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	971	LEU
1	D	864	SER
1	D	694	GLN
1	D	754	ASN
1	D	936	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 39 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	694	GLN
1	D	868	ASN
1	D	728	GLN
1	D	820	ASN
1	D	946	HIS

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](#)

16 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	$\# Z > 2$
3	SO4	A	1102	-	4,4,4	0.34	0	6,6,6	0.12	0
3	SO4	B	1102	-	4,4,4	0.36	0	6,6,6	0.13	0
3	SO4	C	1103	-	4,4,4	0.34	0	6,6,6	0.07	0
3	SO4	D	1104	-	4,4,4	0.34	0	6,6,6	0.12	0
3	SO4	C	1104	-	4,4,4	0.33	0	6,6,6	0.06	0
2	L1S	A	1101	-	40,40,40	0.45	0	56,57,57	1.39	4 (7%)
3	SO4	D	1102	-	4,4,4	0.32	0	6,6,6	0.14	0
2	L1S	B	1101	-	40,40,40	0.43	0	56,57,57	1.35	4 (7%)
3	SO4	A	1103	-	4,4,4	0.33	0	6,6,6	0.09	0
3	SO4	B	1103	-	4,4,4	0.31	0	6,6,6	0.14	0
3	SO4	B	1104	-	4,4,4	0.33	0	6,6,6	0.10	0
3	SO4	D	1103	-	4,4,4	0.32	0	6,6,6	0.12	0
2	L1S	C	1101	-	40,40,40	0.44	0	56,57,57	1.42	4 (7%)
2	L1S	D	1101	-	40,40,40	0.45	0	56,57,57	1.34	4 (7%)
3	SO4	C	1102	-	4,4,4	0.32	0	6,6,6	0.13	0
3	SO4	A	1104	-	4,4,4	0.34	0	6,6,6	0.08	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	L1S	D	1101	-	-	3/26/36/36	0/5/5/5
2	L1S	C	1101	-	-	2/26/36/36	0/5/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	L1S	B	1101	-	-	7/26/36/36	0/5/5/5
2	L1S	A	1101	-	-	6/26/36/36	0/5/5/5

There are no bond length outliers.

The worst 5 of 16 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1101	L1S	C13-C19-N5	-6.69	105.34	109.91
2	D	1101	L1S	C13-C19-N5	-6.29	105.61	109.91
2	B	1101	L1S	C13-C19-N5	-6.24	105.64	109.91
2	C	1101	L1S	C13-C19-N5	-5.90	105.88	109.91
2	A	1101	L1S	C19-C13-N3	5.48	109.97	105.40

There are no chirality outliers.

5 of 18 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1101	L1S	C2-C1-O-C
2	B	1101	L1S	C2-C1-O-C
2	A	1101	L1S	O1-C1-O-C
2	B	1101	L1S	O1-C1-O-C
2	D	1101	L1S	C2-C1-O-C

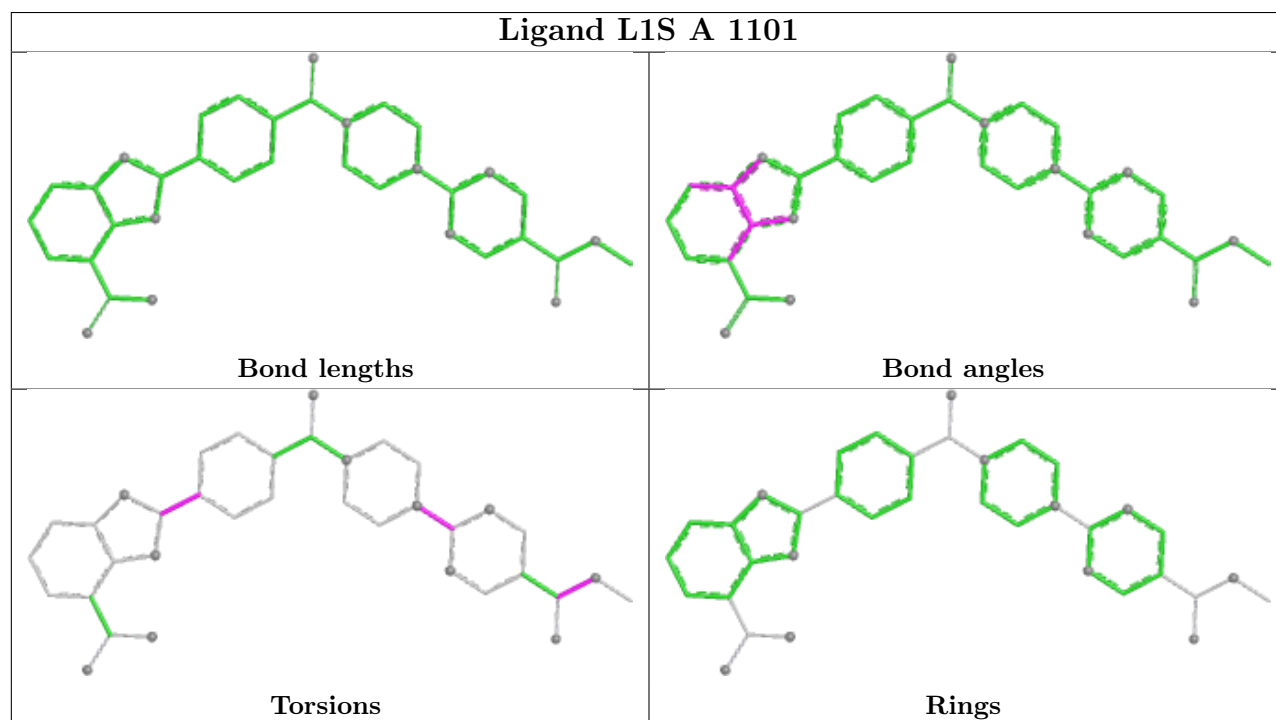
There are no ring outliers.

9 monomers are involved in 14 short contacts:

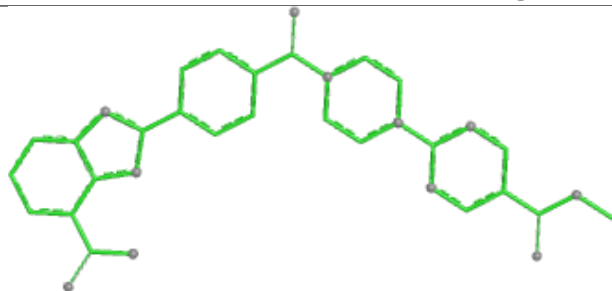
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1103	SO4	1	0
3	D	1104	SO4	1	0
3	C	1104	SO4	1	0
3	B	1103	SO4	1	0
3	B	1104	SO4	1	0
3	D	1103	SO4	1	0
2	C	1101	L1S	3	0
2	D	1101	L1S	4	0
3	A	1104	SO4	1	0

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

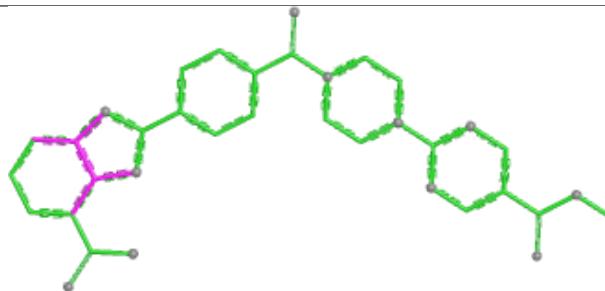
also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted 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.



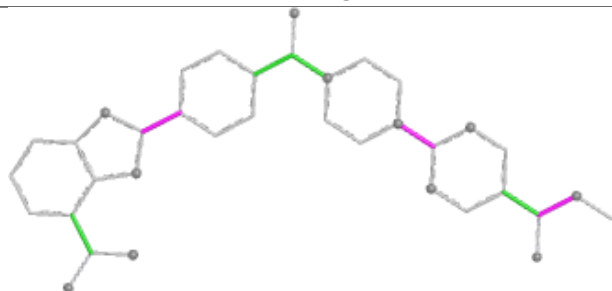
Ligand L1S B 1101



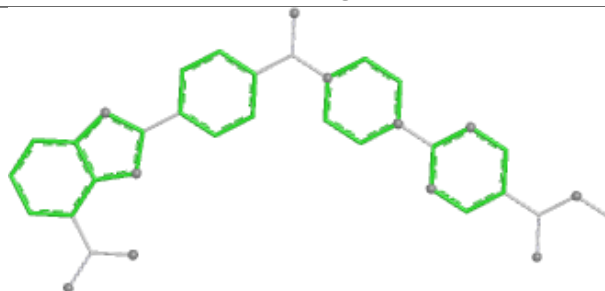
Bond lengths



Bond angles

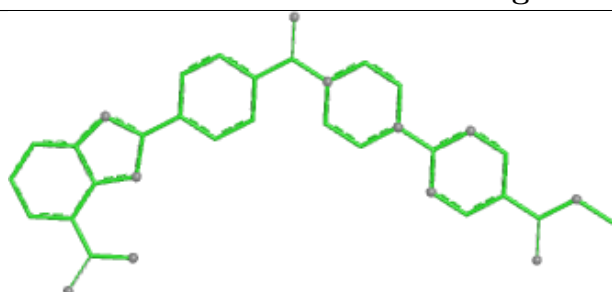


Torsions

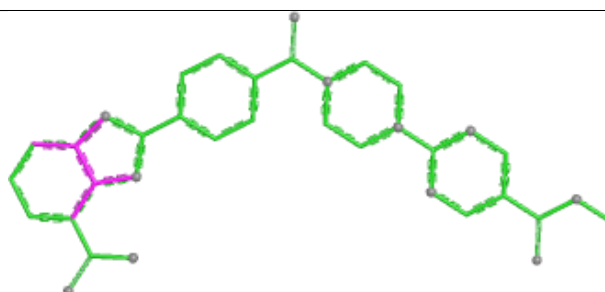


Rings

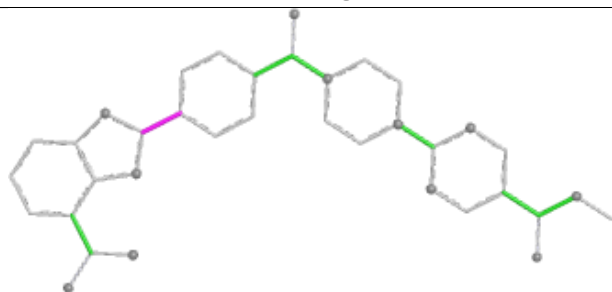
Ligand L1S C 1101



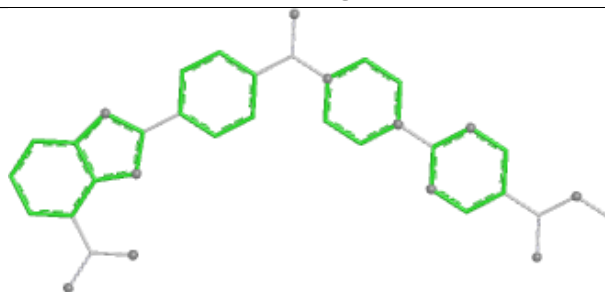
Bond lengths



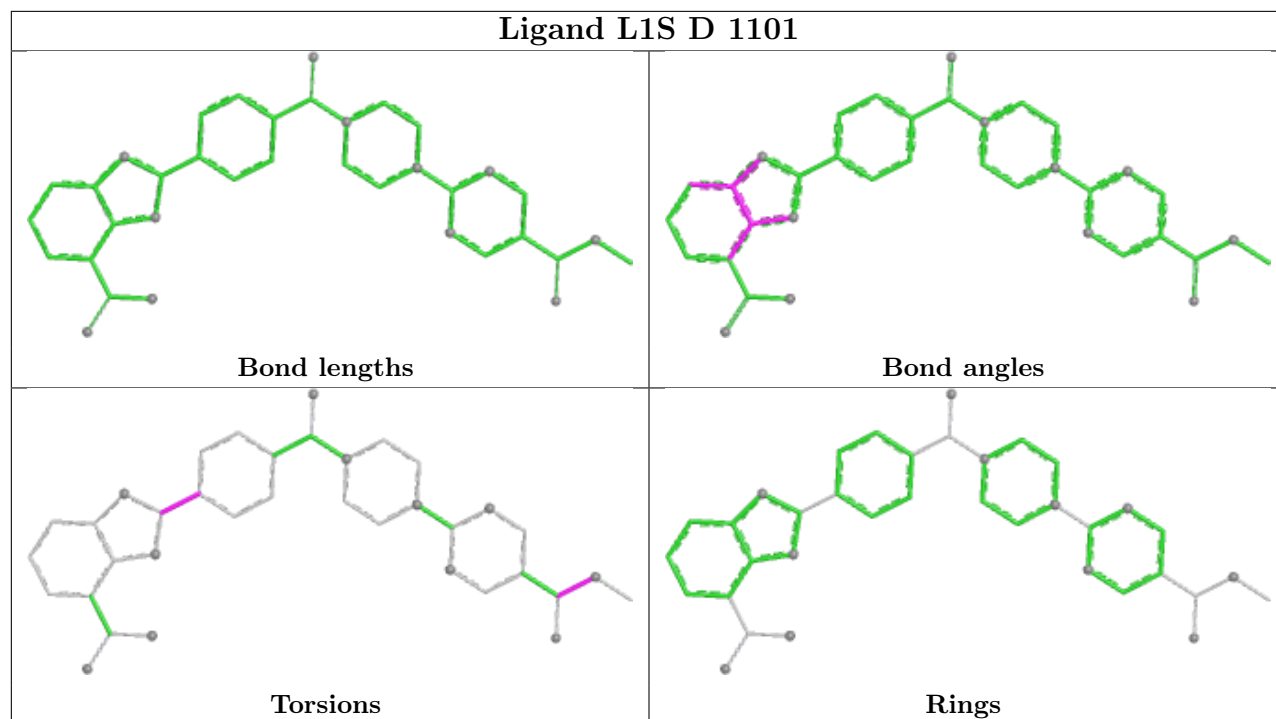
Bond angles



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 [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	340/372 (91%)	1.78	124 (36%) 1 0	34, 97, 161, 180	1 (0%)
1	B	340/372 (91%)	1.78	121 (35%) 1 0	37, 92, 154, 203	1 (0%)
1	C	340/372 (91%)	1.64	113 (33%) 1 1	37, 86, 157, 193	1 (0%)
1	D	340/372 (91%)	1.63	109 (32%) 1 1	39, 91, 160, 194	1 (0%)
All	All	1360/1488 (91%)	1.71	467 (34%) 1 1	34, 92, 159, 203	4 (0%)

The worst 5 of 467 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	888	GLY	7.1
1	B	705	GLN	6.7
1	A	705	GLN	6.2
1	D	705	GLN	6.0
1	C	705	GLN	5.8

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](#)

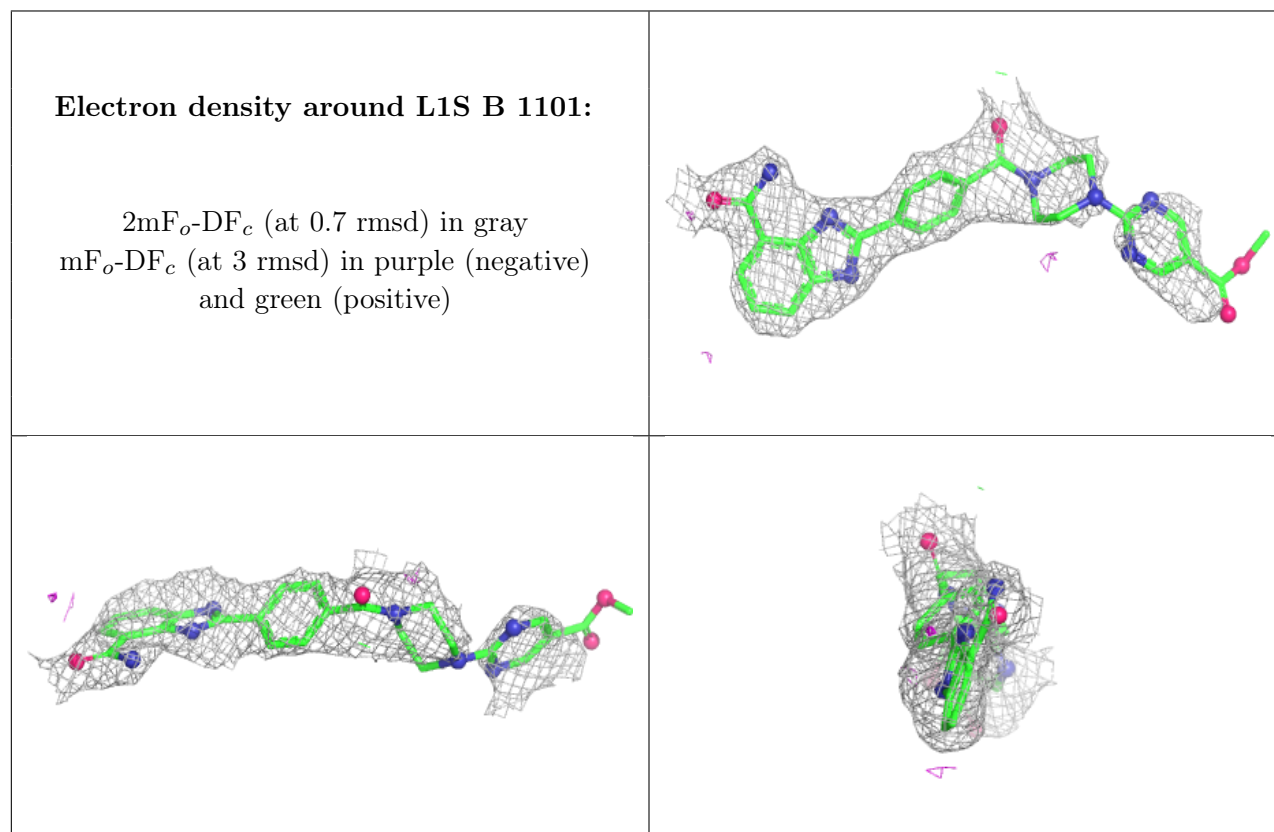
There are no oligosaccharides in this entry.

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.

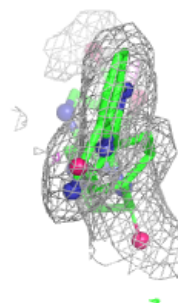
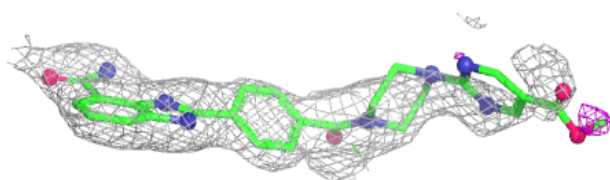
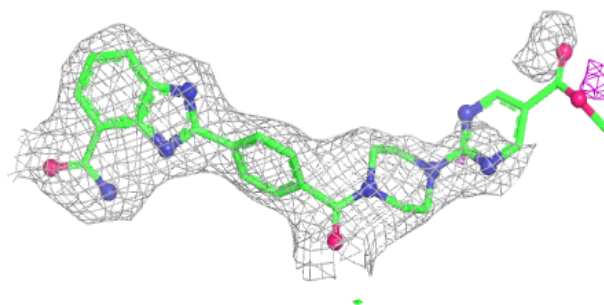
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SO4	A	1104	5/5	0.75	0.16	91,95,99,103	0
3	SO4	D	1104	5/5	0.80	0.14	91,93,95,96	0
2	L1S	B	1101	36/36	0.82	0.17	65,80,141,153	0
3	SO4	B	1104	5/5	0.83	0.15	87,91,95,99	0
3	SO4	C	1104	5/5	0.84	0.16	79,86,87,88	0
2	L1S	A	1101	36/36	0.84	0.15	50,80,145,156	0
3	SO4	A	1103	5/5	0.86	0.18	87,92,97,100	0
3	SO4	B	1103	5/5	0.87	0.13	73,82,88,88	0
3	SO4	D	1103	5/5	0.87	0.18	78,79,81,92	0
2	L1S	D	1101	36/36	0.87	0.13	54,72,122,130	0
3	SO4	B	1102	5/5	0.90	0.16	91,94,96,100	0
2	L1S	C	1101	36/36	0.90	0.12	46,72,131,134	0
3	SO4	C	1103	5/5	0.91	0.16	77,81,84,88	0
3	SO4	C	1102	5/5	0.91	0.14	78,79,82,83	0
3	SO4	D	1102	5/5	0.92	0.17	82,85,88,90	0
3	SO4	A	1102	5/5	0.94	0.11	90,94,99,101	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.

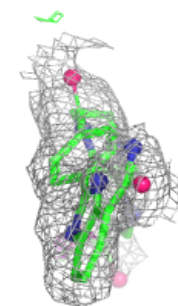
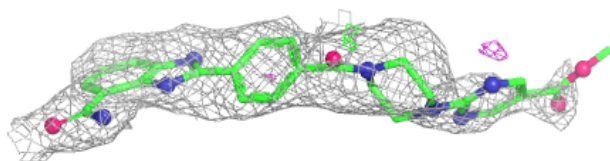
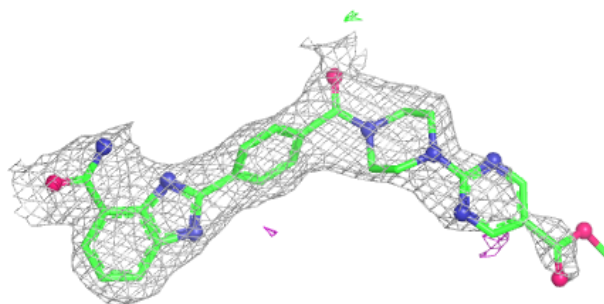


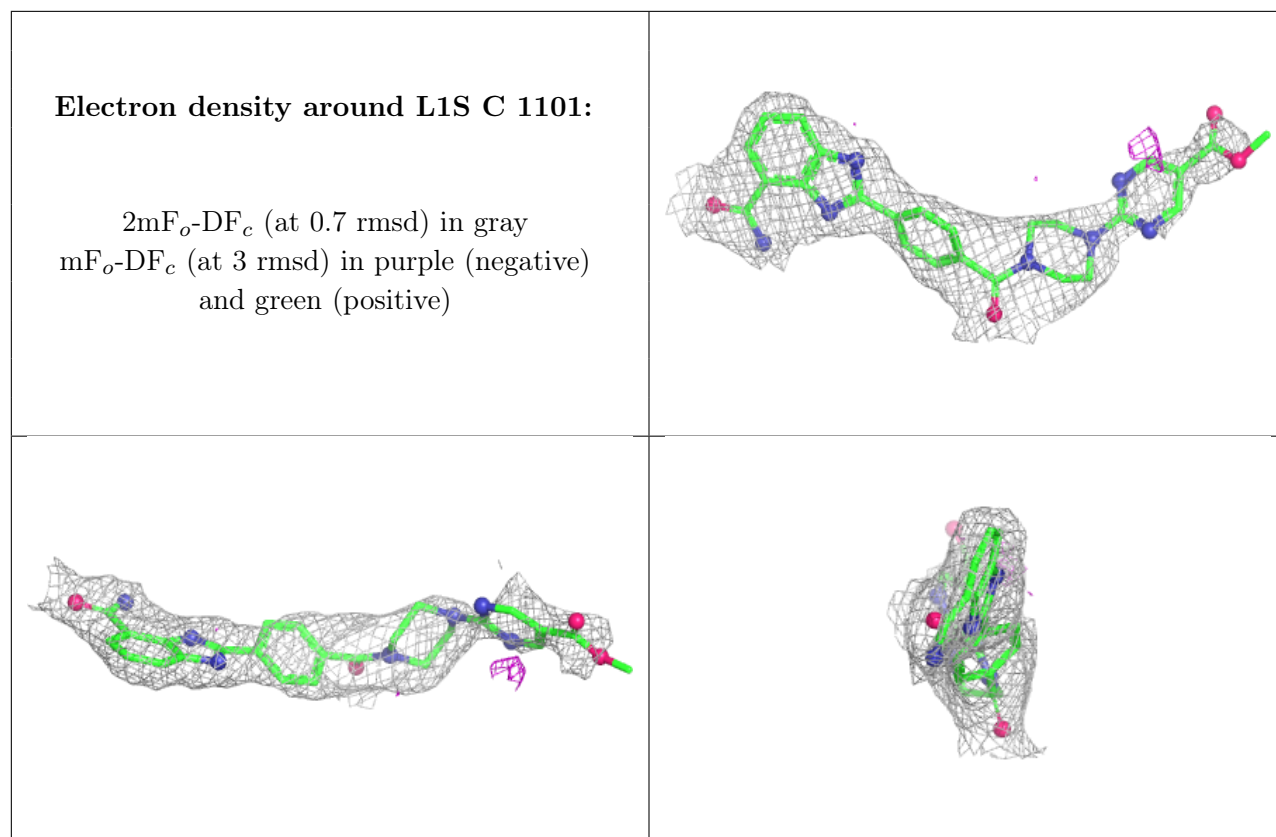
Electron density around L1S A 1101:

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

**Electron density around L1S D 1101:**

$2mF_o-DF_c$ (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.