Full wwPDB X-ray Structure Validation Report ⓘ

Oct 25, 2026 – 10:11 AM EDT

PDB ID	:	9YYF / pdb_00009yyf
Title	:	Crystal structure of M. tuberculosis ProRS bound to compound 4b
Authors	:	Gade, P.; Michalska, K.; Joachimiak, A.; Center for Structural Biology of Infectious Diseases (CSBID)
Deposited on	:	2025-10-28
Resolution	:	2.10 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:

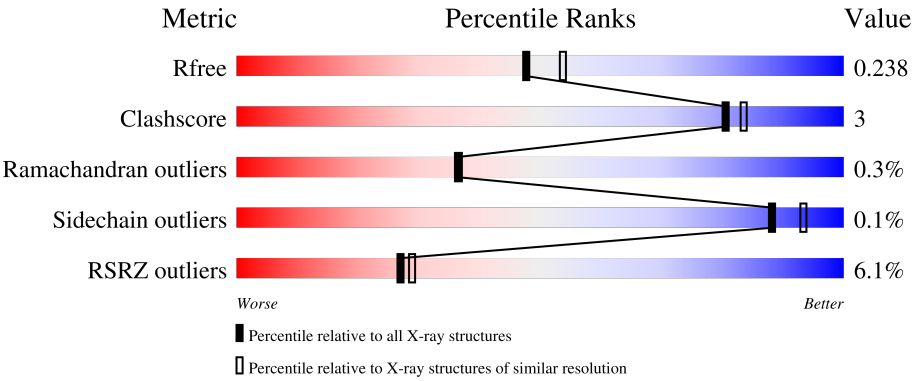
MolProbity	:	4-5-2 with Phenix2.2
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.2
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal 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 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	588	8% 90% 9%
1	B	588	4% 93% 5%

2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 9607 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.

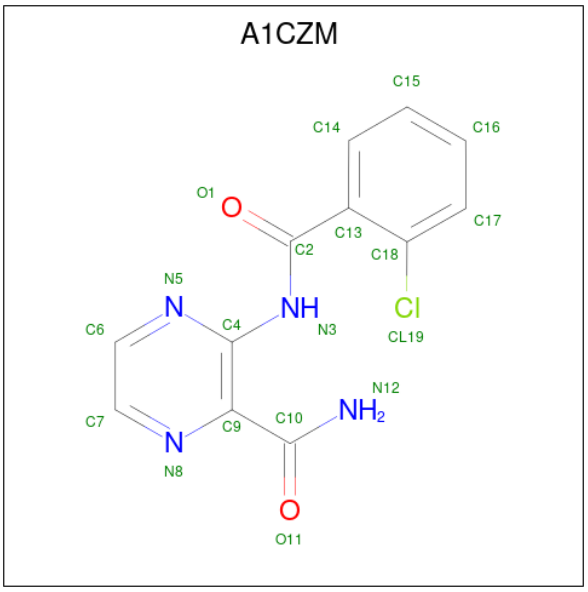
- Molecule 1 is a protein called Proline–tRNA ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	580	Total	C	N	O	S	0	1	0
			4415	2792	778	837	8			
1	B	578	Total	C	N	O	S	0	2	0
			4461	2816	792	845	8			

There are 12 discrepancies between the modelled and reference sequences:

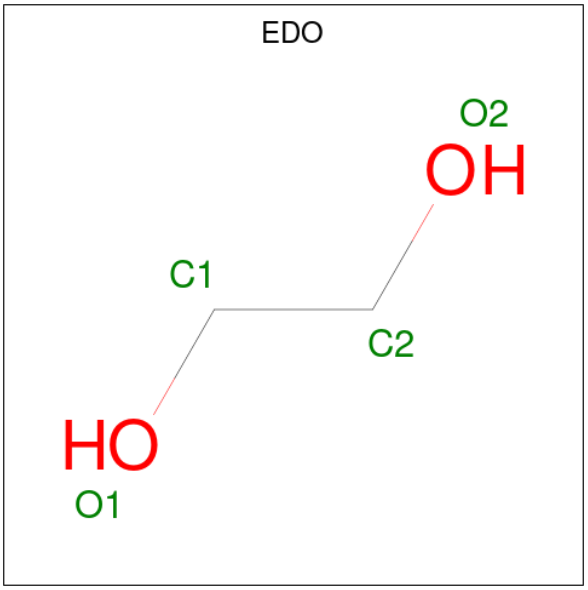
Chain	Residue	Modelled	Actual	Comment	Reference
A	583	GLU	-	expression tag	UNP P9WFT9
A	584	ASN	-	expression tag	UNP P9WFT9
A	585	LEU	-	expression tag	UNP P9WFT9
A	586	TYR	-	expression tag	UNP P9WFT9
A	587	PHE	-	expression tag	UNP P9WFT9
A	588	GLN	-	expression tag	UNP P9WFT9
B	583	GLU	-	expression tag	UNP P9WFT9
B	584	ASN	-	expression tag	UNP P9WFT9
B	585	LEU	-	expression tag	UNP P9WFT9
B	586	TYR	-	expression tag	UNP P9WFT9
B	587	PHE	-	expression tag	UNP P9WFT9
B	588	GLN	-	expression tag	UNP P9WFT9

- Molecule 2 is 3-(2-chlorobenzamido)pyrazine-2-carboxamide (CCD ID: A1CZM) (formula: $C_{12}H_9ClN_4O_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Cl	N	O	0	0
			19	12	1	4	2		
2	B	1	Total	C	Cl	N	O	0	0
			19	12	1	4	2		

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



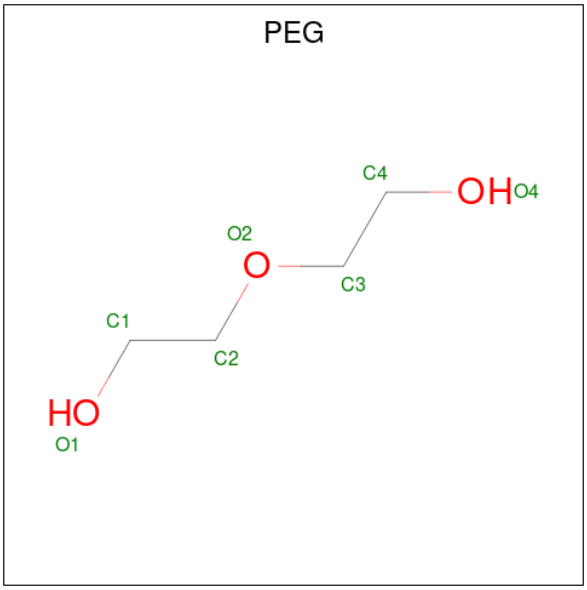
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0

- Molecule 4 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).

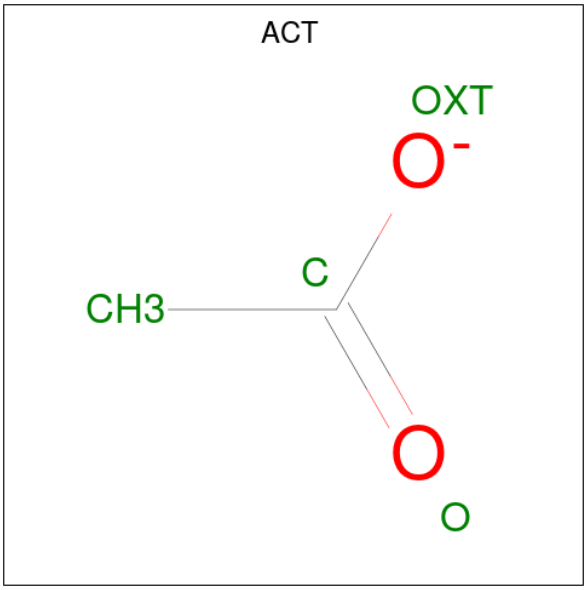


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 7 4 3	0	0

- Molecule 5 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	2	Total Ca 2 2	0	0
5	B	1	Total Ca 1 1	0	0

- Molecule 6 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



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

- Molecule 7 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total Mg 1 1	0	0
7	B	1	Total Mg 1 1	0	0

- Molecule 8 is water.

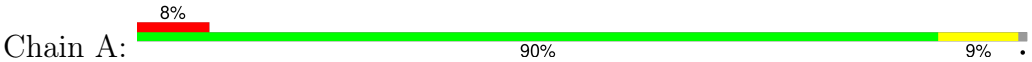
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	322	Total O 322 322	0	0
8	B	315	Total O 315 315	0	0

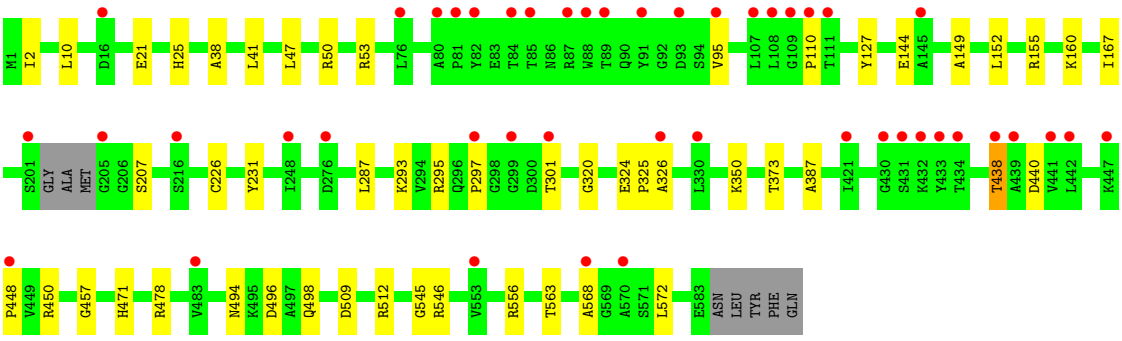
3 Residue-property plots (i)

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: Proline–tRNA ligase

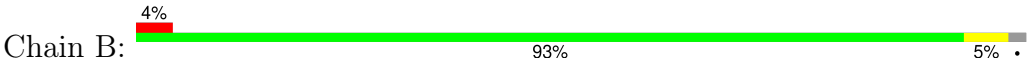
Chain A:

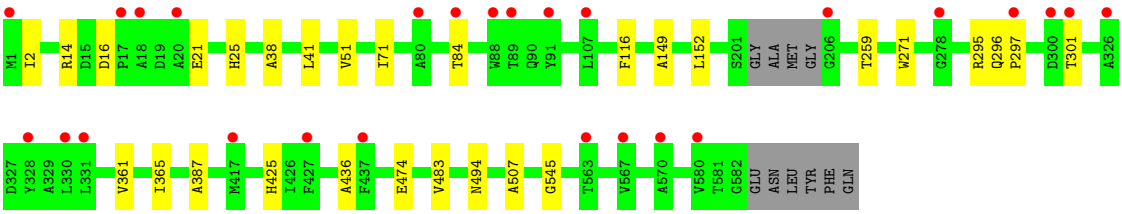




• Molecule 1: Proline–tRNA ligase

Chain B:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	86.02Å 60.12Å 138.65Å 90.00° 90.64° 90.00°	Depositor
Resolution (Å)	49.27 – 2.10 49.27 – 2.10	Depositor EDS
% Data completeness (in resolution range)	98.3 (49.27-2.10) 98.4 (49.27-2.10)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.67 (at 2.10Å)	Xtriage
Refinement program	PHENIX (1.21.2_5419: ???)	Depositor
R, R_{free}	0.198 , 0.238 0.198 , 0.238	Depositor DCC
R_{free} test set	4010 reflections (4.81%)	wwPDB-VP
Wilson B-factor (Å^2)	43.4	Xtriage
Anisotropy	0.265	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 38.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.018 for h,-k,-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	9607	wwPDB-VP
Average B, all atoms (Å^2)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.48% 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: ACT, MG, PEG, A1CZM, CA, EDO

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	0.10	0/4508	0.27	0/6136
1	B	0.10	0/4555	0.26	0/6193
All	All	0.10	0/9063	0.27	0/12329

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	4415	0	4350	31	0
1	B	4461	0	4425	16	0
2	A	19	0	0	0	0
2	B	19	0	0	0	0
3	A	20	0	30	2	0
3	B	20	0	30	2	0
4	A	7	0	10	2	0
5	A	2	0	0	0	0
5	B	1	0	0	0	0
6	A	4	0	3	0	0
7	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	B	1	0	0	0	0
8	A	322	0	0	0	0
8	B	315	0	0	0	0
All	All	9607	0	8848	47	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 3.

All (47) 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:387:ALA:HB1	3:A:606:EDO:H21	1.82	0.62
1:B:84:THR:HG22	1:B:436:ALA:HB1	1.85	0.59
1:A:568:ALA:HA	1:A:572:LEU:HD22	1.85	0.58
1:B:296:GLN:HG3	1:B:297:PRO:HD2	1.86	0.58
1:B:295:ARG:HB2	1:B:301:THR:HG22	1.88	0.55
1:A:127:TYR:HB2	1:A:167:ILE:HG21	1.88	0.54
1:B:474:GLU:H	1:B:474:GLU:CD	2.15	0.53
1:B:14:ARG:HH12	1:B:16:ASP:HB2	1.72	0.52
1:B:38:ALA:HB3	1:B:41:LEU:HD12	1.91	0.52
1:A:50:ARG:HA	4:A:607:PEG:H31	1.93	0.51
1:B:545:GLY:HA3	3:B:606:EDO:H11	1.93	0.50
1:B:494:ASN:HB2	1:B:545:GLY:HA2	1.93	0.50
1:B:387:ALA:HB1	3:B:602:EDO:H12	1.93	0.49
1:A:350:LYS:HD2	3:A:602:EDO:H12	1.93	0.49
1:A:546:ARG:CZ	1:A:546:ARG:H	2.26	0.49
1:A:556:ARG:HG2	1:A:563:THR:HG23	1.94	0.48
1:A:440:ASP:HB3	1:A:448:PRO:HB3	1.95	0.48
1:A:53:ARG:HG3	4:A:607:PEG:H32	1.96	0.48
1:A:38:ALA:HB3	1:A:41:LEU:HD12	1.96	0.48
1:A:320:GLY:O	1:A:324:GLU:HG2	2.15	0.47
1:B:259:THR:HG22	1:B:271:TRP:CE2	2.50	0.47
1:B:21:GLU:HG2	1:B:25:HIS:HD2	1.79	0.47
1:A:295:ARG:HH21	1:A:297:PRO:HA	1.79	0.47
1:A:293:LYS:HG2	1:A:301:THR:HB	1.98	0.45
1:A:509:ASP:HA	1:A:512:ARG:HE	1.81	0.45
1:B:51:VAL:HG13	1:B:483:VAL:HG12	1.99	0.45
1:A:21:GLU:HG2	1:A:25:HIS:HD2	1.82	0.44
1:B:149:ALA:HB1	1:B:152:LEU:HD12	1.97	0.44
1:A:471:HIS:O	1:A:478:ARG:HB2	2.17	0.44
1:A:149:ALA:HB1	1:A:152:LEU:HD12	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:95:VAL:HG11	1:A:110:PRO:HG3	2.00	0.44
1:A:494:ASN:HB2	1:A:545:GLY:HA2	2.01	0.43
1:A:287:LEU:HB3	1:A:373:THR:HB	2.01	0.43
1:A:496:ASP:OD2	1:A:498:GLN:HG2	2.18	0.43
1:A:295:ARG:HB3	1:A:301:THR:HG22	2.01	0.43
1:A:160:LYS:HE3	1:A:160:LYS:HB3	1.90	0.43
1:A:160:LYS:O	1:A:457:GLY:HA2	2.19	0.43
1:B:361:VAL:HB	1:B:365:ILE:HG13	2.01	0.42
1:A:324:GLU:HA	1:A:326:ALA:H	1.85	0.42
1:A:226:CYS:HB3	1:A:231:TYR:HB3	2.01	0.41
1:A:509:ASP:O	1:A:512:ARG:HG2	2.19	0.41
1:A:10:LEU:HD22	1:A:47:LEU:HD11	2.03	0.41
1:A:438:THR:HA	1:A:450:ARG:NH2	2.36	0.41
1:B:2:ILE:HD11	1:B:507:ALA:O	2.21	0.41
1:A:2:ILE:HD13	1:A:2:ILE:HA	1.97	0.40
1:A:144:GLU:H	1:A:155:ARG:HG2	1.86	0.40
1:B:71:ILE:HD11	1:B:116:PHE:CE2	2.56	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	577/588 (98%)	552 (96%)	22 (4%)	3 (0%)	24 22
1	B	576/588 (98%)	559 (97%)	17 (3%)	0	100 100
All	All	1153/1176 (98%)	1111 (96%)	39 (3%)	3 (0%)	36 36

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	207	SER

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Mol	Chain	Res	Type
1	A	438	THR
1	A	325	PRO

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	448/465 (96%)	448 (100%)	0	100	100
1	B	459/465 (99%)	457 (100%)	2 (0%)	84	89
All	All	907/930 (98%)	905 (100%)	2 (0%)	88	93

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	425[A]	HIS
1	B	425[B]	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	HIS
1	A	562	GLN
1	B	489	HIS

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)

Of 19 ligands modelled in this entry, 5 are monoatomic - leaving 14 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	$\#\|Z\| > 2$
4	PEG	A	607	-	6,6,6	0.26	0	5,5,5	0.26	0
3	EDO	A	605	-	3,3,3	0.24	0	2,2,2	0.34	0
3	EDO	A	603	-	3,3,3	0.27	0	2,2,2	0.27	0
3	EDO	A	606	-	3,3,3	0.25	0	2,2,2	0.35	0
3	EDO	B	602	-	3,3,3	0.25	0	2,2,2	0.35	0
3	EDO	B	605	-	3,3,3	0.25	0	2,2,2	0.33	0
2	A1CZM	A	601	-	20,20,20	0.15	0	23,27,27	0.34	0
3	EDO	B	606	-	3,3,3	0.25	0	2,2,2	0.33	0
2	A1CZM	B	601	-	20,20,20	0.13	0	23,27,27	0.33	0
3	EDO	B	603	-	3,3,3	0.24	0	2,2,2	0.31	0
3	EDO	A	604	-	3,3,3	0.25	0	2,2,2	0.32	0
3	EDO	B	604	-	3,3,3	0.24	0	2,2,2	0.39	0
6	ACT	A	610	-	3,3,3	1.13	0	3,3,3	1.22	0
3	EDO	A	602	-	3,3,3	0.24	0	2,2,2	0.32	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
4	PEG	A	607	-	-	2/4/4/4	-
3	EDO	A	605	-	-	1/1/1/1	-
3	EDO	A	603	-	-	1/1/1/1	-
3	EDO	A	606	-	-	0/1/1/1	-
3	EDO	B	602	-	-	0/1/1/1	-
3	EDO	B	605	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	A1CZM	A	601	-	-	0/12/12/12	0/2/2/2
3	EDO	B	606	-	-	0/1/1/1	-
2	A1CZM	B	601	-	-	0/12/12/12	0/2/2/2
3	EDO	B	603	-	-	1/1/1/1	-
3	EDO	A	604	-	-	0/1/1/1	-
3	EDO	B	604	-	-	0/1/1/1	-
3	EDO	A	602	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	607	PEG	O1-C1-C2-O2
3	A	603	EDO	O1-C1-C2-O2
3	B	603	EDO	O1-C1-C2-O2
3	A	605	EDO	O1-C1-C2-O2
4	A	607	PEG	C1-C2-O2-C3

There are no ring outliers.

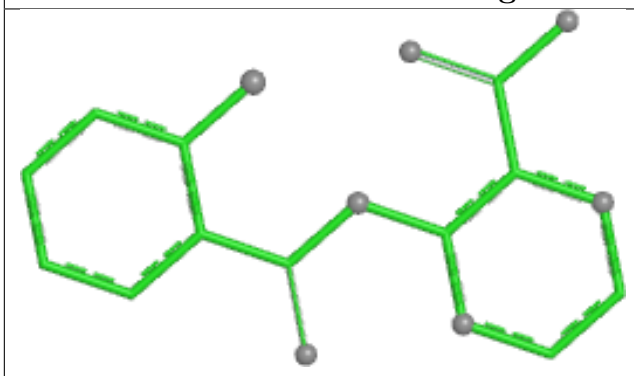
5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	607	PEG	2	0
3	A	606	EDO	1	0
3	B	602	EDO	1	0
3	B	606	EDO	1	0
3	A	602	EDO	1	0

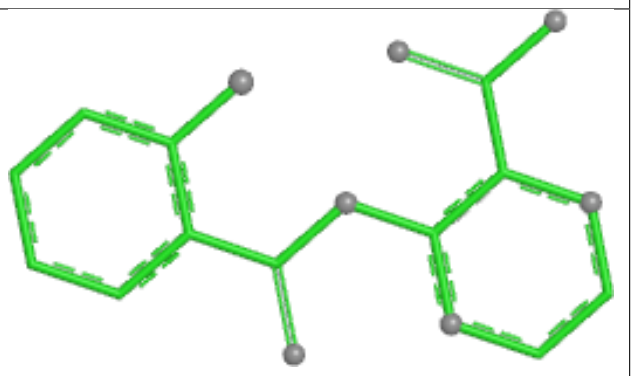
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

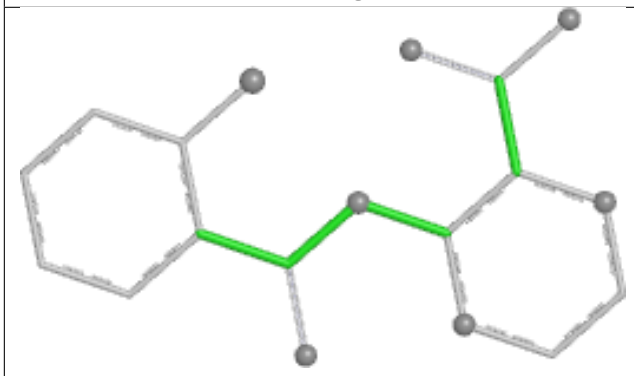
Ligand A1CZM A 601



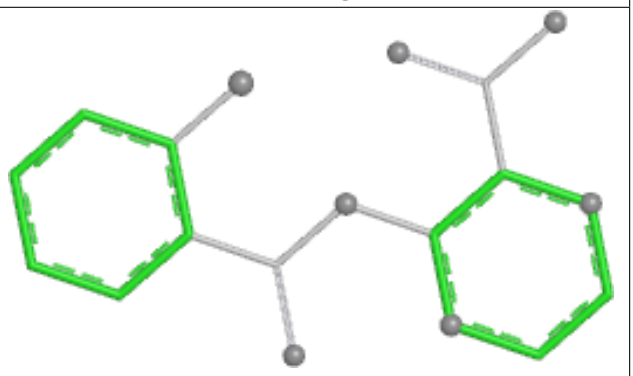
Bond lengths



Bond angles

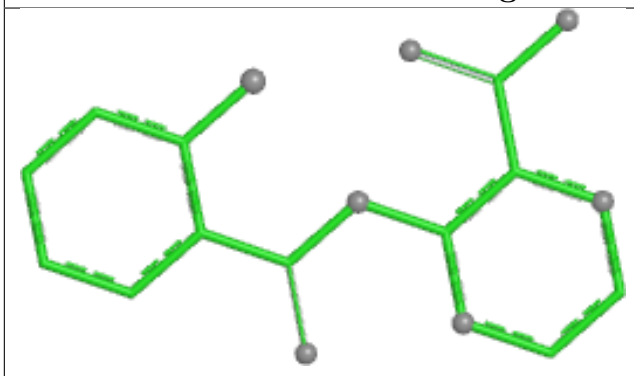


Torsions

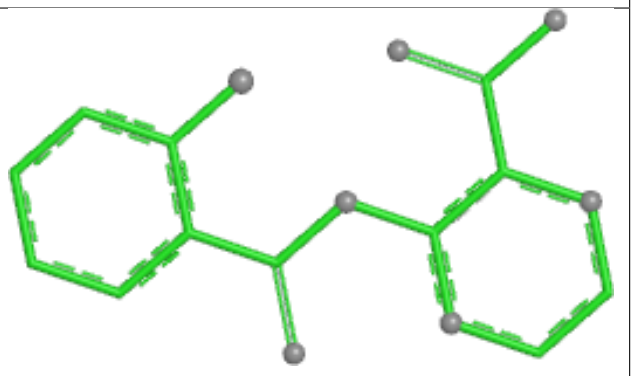


Rings

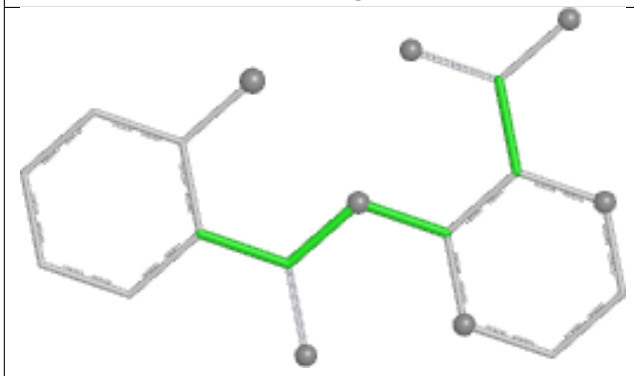
Ligand A1CZM B 601



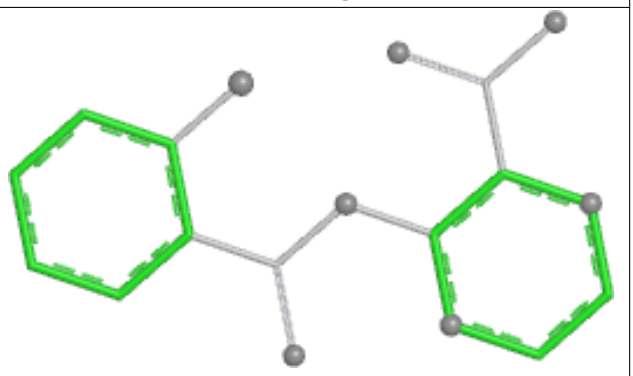
Bond lengths



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, 95^{th} 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	580/588 (98%)	0.42	45 (7%)	19	20	26, 49, 99, 123	2 (0%)
1	B	578/588 (98%)	0.33	26 (4%)	38	40	24, 50, 88, 110	2 (0%)
All	All	1158/1176 (98%)	0.37	71 (6%)	27	29	24, 50, 94, 123	4 (0%)

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	205	GLY	4.6
1	A	248	ILE	4.4
1	A	82	TYR	3.9
1	A	111	THR	3.9
1	A	81	PRO	3.6
1	A	448	PRO	3.6
1	A	431	SER	3.5
1	A	276	ASP	3.5
1	A	110	PRO	3.5
1	A	434	THR	3.5
1	A	109	GLY	3.4
1	A	433	TYR	3.3
1	A	88	TRP	3.2
1	B	84	THR	3.2
1	A	93	ASP	3.1
1	A	430	GLY	3.0
1	B	300	ASP	3.0
1	B	107	LEU	3.0
1	A	87	ARG	2.9
1	A	301	THR	2.9
1	A	330	LEU	2.8
1	B	570	ALA	2.8
1	A	91	TYR	2.7
1	A	107	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	80	ALA	2.7
1	B	89	THR	2.7
1	A	568	ALA	2.6
1	A	85	THR	2.6
1	A	439	ALA	2.6
1	A	421	ILE	2.6
1	A	442	LEU	2.6
1	A	89	THR	2.5
1	B	1	MET	2.5
1	A	201	SER	2.5
1	A	432	LYS	2.5
1	A	438	THR	2.4
1	B	297	PRO	2.4
1	B	20	ALA	2.4
1	A	76	LEU	2.4
1	B	437	PHE	2.4
1	A	216	SER	2.4
1	B	331	LEU	2.4
1	A	326	ALA	2.3
1	A	80	ALA	2.3
1	A	299	GLY	2.3
1	A	95	VAL	2.3
1	A	441	VAL	2.3
1	B	17	PRO	2.3
1	B	563	THR	2.3
1	B	427	PHE	2.2
1	A	483	VAL	2.2
1	B	580	VAL	2.2
1	A	570	ALA	2.2
1	B	417	MET	2.1
1	B	567	VAL	2.1
1	A	108	LEU	2.1
1	A	447	LYS	2.1
1	A	84	THR	2.1
1	A	297	PRO	2.1
1	B	206	GLY	2.1
1	A	553	VAL	2.1
1	B	326	ALA	2.1
1	B	91	TYR	2.1
1	B	301	THR	2.1
1	B	328	TYR	2.1
1	B	278	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	145	ALA	2.1
1	B	18	ALA	2.0
1	B	88	TRP	2.0
1	B	330	LEU	2.0
1	A	16	ASP	2.0

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, 95^{th} 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
6	ACT	A	610	4/4	0.69	0.14	74,75,83,85	0
3	EDO	B	603	4/4	0.71	0.20	56,61,66,75	0
4	PEG	A	607	7/7	0.80	0.17	48,51,65,69	0
3	EDO	A	604	4/4	0.81	0.12	66,68,73,78	0
3	EDO	B	605	4/4	0.82	0.09	75,79,80,83	0
7	MG	A	611	1/1	0.84	0.10	76,76,76,76	0
3	EDO	A	603	4/4	0.85	0.21	53,60,67,75	0
3	EDO	A	602	4/4	0.86	0.17	56,56,58,67	0
5	CA	B	607	1/1	0.88	0.09	91,91,91,91	0
3	EDO	B	606	4/4	0.88	0.12	67,72,75,75	0
3	EDO	B	604	4/4	0.88	0.13	62,63,66,67	0
7	MG	B	608	1/1	0.89	0.06	81,81,81,81	0
3	EDO	B	602	4/4	0.91	0.13	54,54,57,57	0
2	A1CZM	A	601	19/19	0.92	0.09	45,50,62,66	0
3	EDO	A	605	4/4	0.93	0.12	50,51,61,74	0
5	CA	A	608	1/1	0.93	0.07	94,94,94,94	0
2	A1CZM	B	601	19/19	0.93	0.09	35,48,64,66	0
3	EDO	A	606	4/4	0.94	0.10	46,49,55,56	0

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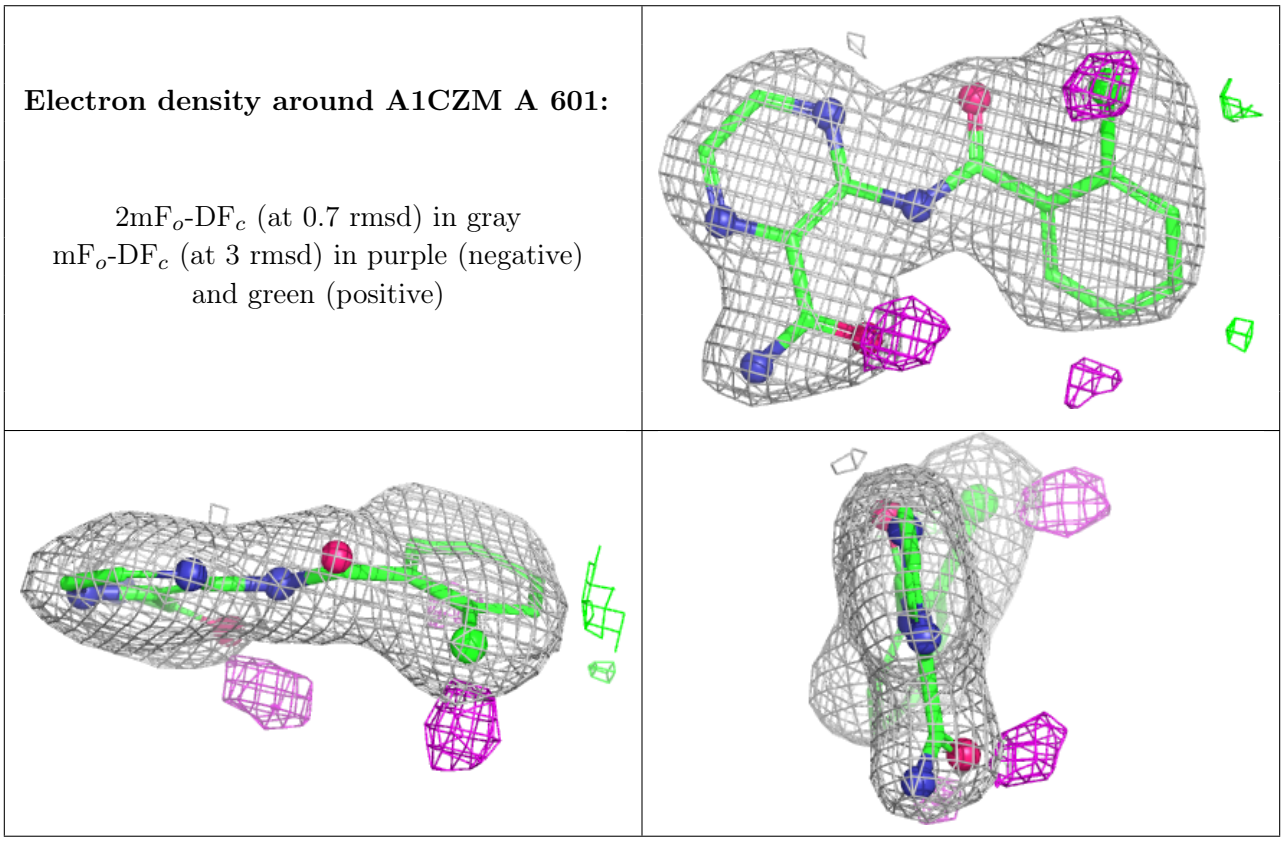
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	CA	A	609	1/1	0.97	0.04	78,78,78,78	0

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.

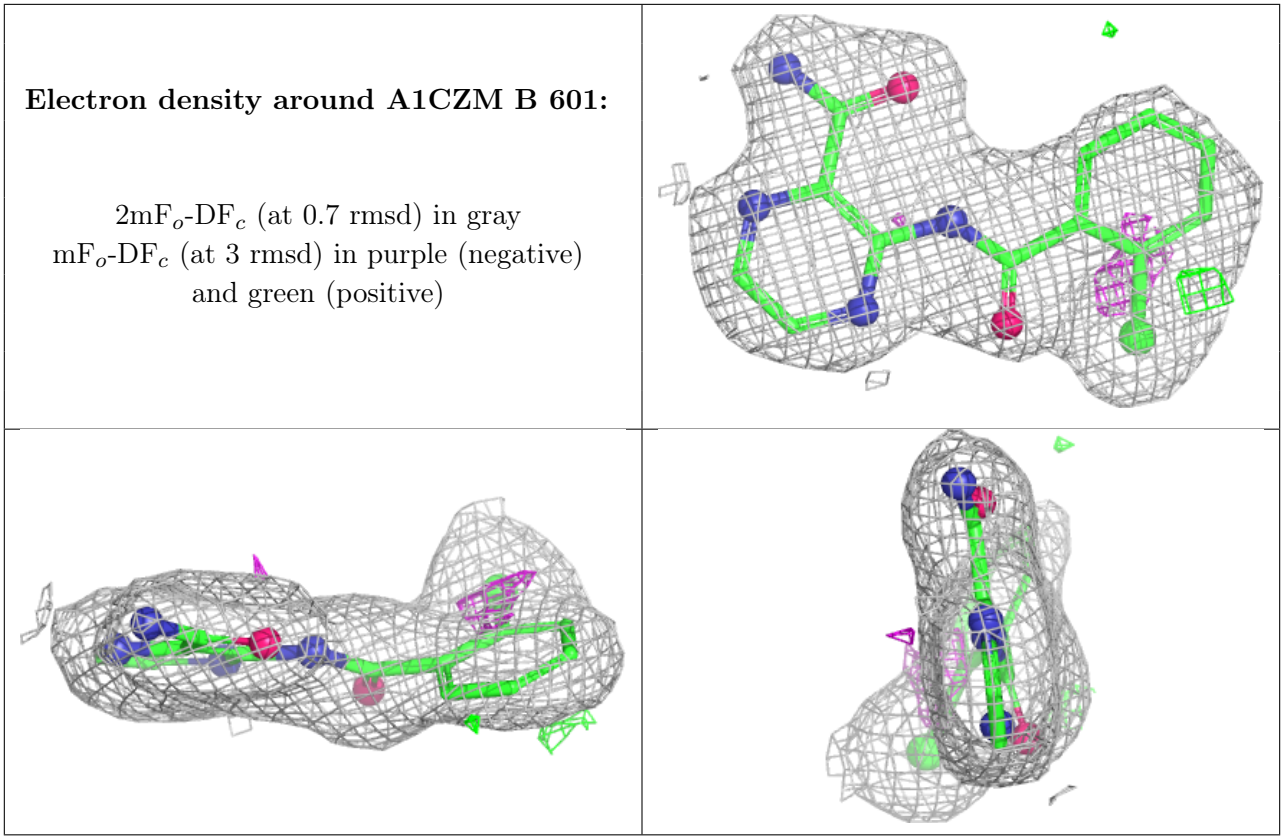
Electron density around A1CZM A 601:

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



Electron density around A1CZM B 601:

$2mF_o$-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.