



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 07:15 PM UTC

PDB ID : 2B51 / pdb_00002b51
Title : Structural Basis for UTP Specificity of RNA Editing TUTases from Trypanosoma Brucei
Authors : Deng, J.; Ernst, N.L.; Turley, S.; Stuart, K.D.; Hol, W.G.
Deposited on : 2005-09-27
Resolution : 2.05 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

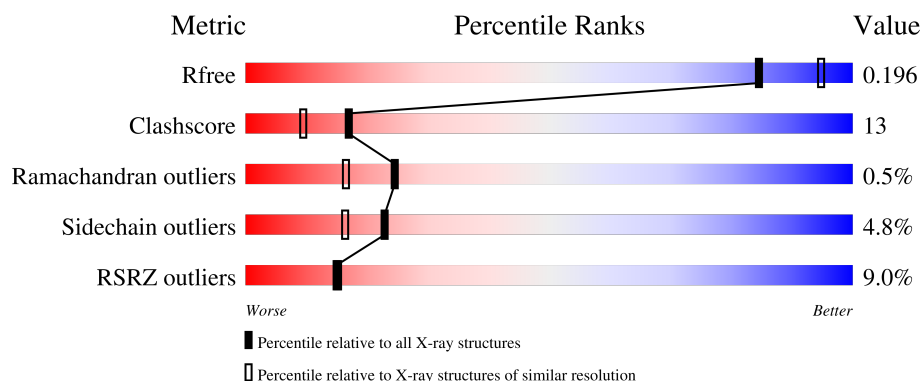
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	468	9% 76% 16% • 5%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4000 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 RNA editing complex protein MP57.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	444	Total	C	N	O	S	11	1	0
			3587	2277	642	644	24			

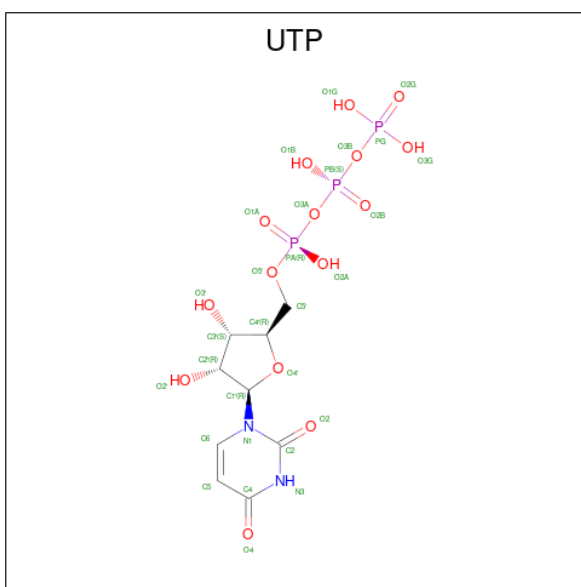
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	94	ARG	TRP	engineered mutation	UNP Q86MV5
A	122	ARG	TRP	engineered mutation	UNP Q86MV5
A	207	ARG	TRP	engineered mutation	UNP Q86MV5

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

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

- Molecule 3 is URIDINE 5'-TRIPHOSPHATE (CCD ID: UTP) (formula: C₉H₁₅N₂O₁₅P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 29	C 9	N 2	O 15	P 3	0	0
3	A	1	Total 29	C 9	N 2	O 15	P 3	0	0
3	A	1	Total 29	C 9	N 2	O 15	P 3	0	0

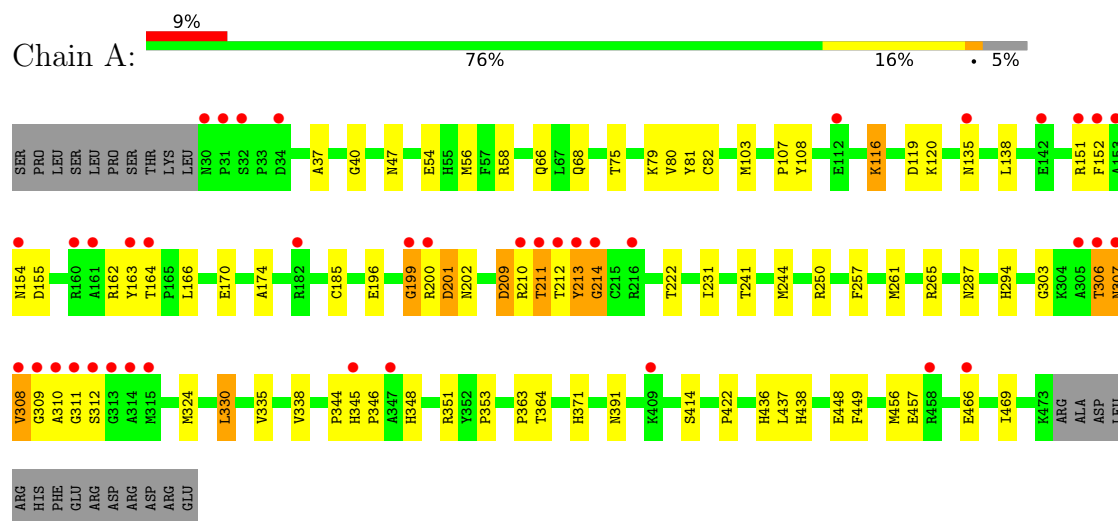
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	325	Total O 325 325	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: RNA editing complex protein MP57



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	91.01Å 91.01Å 162.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.05 50.00 – 2.05	Depositor EDS
% Data completeness (in resolution range)	99.9 (50.00-2.05) 99.9 (50.00-2.05)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.187 , 0.227 0.192 , 0.196	Depositor DCC
R_{free} test set	2313 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	30.3	Xtriage
Anisotropy	0.255	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 42.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4000	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.75% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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	0.64	0/3685	0.83	0/4993

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	5

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	199	GLY	Peptide
1	A	213	TYR	Peptide
1	A	214	GLY	Peptide
1	A	307	ASN	Peptide
1	A	308	VAL	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	3587	0	3487	90	0
2	A	1	0	0	0	0
3	A	87	0	33	4	0
4	A	325	0	0	2	0
All	All	4000	0	3520	91	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 13.

All (91) 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:213:TYR:CD2	1:A:214:GLY:HA2	1.96	1.00
1:A:308:VAL:N	1:A:309:GLY:HA3	1.85	0.92
1:A:211:THR:HG21	1:A:213:TYR:CE1	2.05	0.91
1:A:241:THR:H	1:A:244:MET:HE3	1.42	0.83
1:A:306:THR:CG2	1:A:307:ASN:N	2.41	0.83
1:A:306:THR:HG23	1:A:307:ASN:H	1.44	0.83
1:A:303:GLY:O	1:A:309:GLY:HA2	1.78	0.82
1:A:211:THR:OG1	1:A:213:TYR:CZ	2.34	0.80
1:A:211:THR:CB	1:A:213:TYR:CE1	2.64	0.80
1:A:211:THR:OG1	1:A:214:GLY:HA3	1.85	0.76
1:A:68:GLN:HE22	1:A:80:VAL:H	1.35	0.75
1:A:47:ASN:HD22	1:A:294:HIS:HD2	1.34	0.73
1:A:306:THR:CG2	1:A:307:ASN:H	1.99	0.73
1:A:211:THR:CG2	1:A:213:TYR:CE1	2.71	0.72
1:A:391:ASN:HD21	1:A:422:PRO:HA	1.55	0.70
1:A:201:ASP:O	1:A:222:THR:HG21	1.91	0.69
1:A:414:SER:OG	1:A:438:HIS:HD2	1.75	0.69
1:A:306:THR:HG22	1:A:307:ASN:N	2.08	0.69
1:A:306:THR:HG22	1:A:308:VAL:HG23	1.73	0.69
1:A:241:THR:N	1:A:244:MET:HE3	2.07	0.68
1:A:324:MET:HE1	1:A:449:PHE:CZ	2.29	0.67
1:A:324:MET:HE1	1:A:449:PHE:HZ	1.61	0.66
1:A:47:ASN:HD22	1:A:294:HIS:CD2	2.13	0.65
1:A:308:VAL:N	1:A:309:GLY:CA	2.58	0.65
1:A:211:THR:HG21	1:A:213:TYR:CZ	2.32	0.65
1:A:211:THR:OG1	1:A:214:GLY:CA	2.45	0.64
1:A:371:HIS:HD2	4:A:581:HOH:O	1.81	0.62
1:A:68:GLN:NE2	1:A:80:VAL:H	1.98	0.61
1:A:306:THR:CG2	1:A:308:VAL:HG23	2.31	0.61
1:A:211:THR:OG1	1:A:213:TYR:CE2	2.49	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:196:GLU:OE1	1:A:200:ARG:HD2	2.00	0.60
1:A:199:GLY:HA2	1:A:200:ARG:CG	2.32	0.60
1:A:199:GLY:O	1:A:202:ASN:HB2	2.02	0.59
1:A:306:THR:HG23	1:A:448:GLU:CD	2.28	0.59
1:A:211:THR:HB	1:A:213:TYR:CD1	2.37	0.59
1:A:371:HIS:HE1	1:A:457:GLU:OE2	1.87	0.58
1:A:56:MET:HE1	1:A:163:TYR:CE2	2.39	0.56
1:A:211:THR:OG1	1:A:213:TYR:CE1	2.57	0.55
1:A:211:THR:HB	1:A:213:TYR:CE1	2.42	0.55
1:A:312:SER:O	3:A:502:UTP:H1'	2.06	0.55
1:A:391:ASN:ND2	4:A:517:HOH:O	2.40	0.55
1:A:241:THR:HG23	1:A:244:MET:CE	2.37	0.55
1:A:199:GLY:HA2	1:A:200:ARG:HG2	1.88	0.54
1:A:211:THR:HG21	1:A:213:TYR:HE1	1.68	0.54
1:A:330:LEU:HD13	1:A:335:VAL:HG13	1.90	0.53
1:A:456:MET:HG3	1:A:469:ILE:HD13	1.92	0.52
1:A:213:TYR:CE2	1:A:214:GLY:HA2	2.42	0.52
1:A:308:VAL:H	1:A:309:GLY:HA3	1.68	0.51
1:A:345:HIS:HB3	1:A:348:HIS:HD2	1.75	0.51
1:A:56:MET:HE2	1:A:162:ARG:HB3	1.91	0.51
1:A:213:TYR:CD2	1:A:214:GLY:CA	2.84	0.51
1:A:82:CYS:HB3	1:A:353:PRO:HD3	1.91	0.51
1:A:56:MET:HE1	1:A:163:TYR:HE2	1.75	0.50
1:A:54:GLU:OE2	1:A:58:ARG:NH2	2.34	0.50
1:A:211:THR:CB	1:A:213:TYR:CD1	2.95	0.50
1:A:166:LEU:HB3	1:A:170:GLU:HB2	1.92	0.49
1:A:306:THR:HG23	1:A:448:GLU:OE2	2.12	0.49
1:A:199:GLY:HA2	1:A:200:ARG:CB	2.42	0.49
1:A:344:PRO:HB2	1:A:348:HIS:HB2	1.95	0.49
1:A:66:GLN:NE2	1:A:261:MET:H	2.10	0.48
1:A:231:ILE:HG23	1:A:250:ARG:HB3	1.95	0.48
1:A:56:MET:CE	1:A:163:TYR:CD2	2.96	0.47
1:A:287:ASN:ND2	3:A:503:UTP:O1G	2.45	0.47
1:A:185:CYS:SG	1:A:210:ARG:HG2	2.55	0.46
1:A:56:MET:HE3	1:A:163:TYR:HD2	1.81	0.45
1:A:81:TYR:CD2	1:A:103:MET:HE3	2.51	0.45
1:A:56:MET:CE	1:A:163:TYR:HD2	2.30	0.45
1:A:306:THR:C	1:A:308:VAL:H	2.24	0.45
1:A:241:THR:HG23	1:A:244:MET:HE3	1.99	0.45
1:A:152:PHE:CZ	1:A:154:ASN:HA	2.52	0.44
1:A:107:PRO:HG3	1:A:345:HIS:CE1	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:GLU:O	1:A:58:ARG:HG2	2.17	0.44
1:A:199:GLY:HA2	1:A:200:ARG:HB2	2.00	0.44
1:A:212:THR:HG22	1:A:212:THR:O	2.18	0.43
1:A:211:THR:CG2	1:A:213:TYR:CZ	2.98	0.43
1:A:436:HIS:CE1	3:A:502:UTP:H3'	2.54	0.43
1:A:209:ASP:OD1	1:A:209:ASP:N	2.52	0.42
1:A:138:LEU:HD11	1:A:151:ARG:HD3	2.02	0.42
1:A:116:LYS:HE2	1:A:116:LYS:HB2	1.58	0.42
1:A:37:ALA:HB2	1:A:364:THR:OG1	2.20	0.42
1:A:345:HIS:HB3	1:A:348:HIS:CD2	2.55	0.42
1:A:456:MET:CG	1:A:469:ILE:HD13	2.50	0.41
1:A:199:GLY:CA	1:A:200:ARG:HB2	2.50	0.41
1:A:66:GLN:HE22	1:A:261:MET:H	1.69	0.41
1:A:311:GLY:HA3	1:A:312:SER:HA	1.77	0.40
1:A:107:PRO:HG2	1:A:108:TYR:CE2	2.56	0.40
1:A:174:ALA:HA	1:A:257:PHE:O	2.21	0.40
3:A:502:UTP:O5'	3:A:502:UTP:H6	2.04	0.40
1:A:414:SER:OG	1:A:438:HIS:CD2	2.64	0.40
1:A:40:GLY:HA3	1:A:363:PRO:O	2.20	0.40
1:A:308:VAL:H	1:A:309:GLY:CA	2.29	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	443/468 (95%)	427 (96%)	14 (3%)	2 (0%)	24 16

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	201	ASP

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Mol	Chain	Res	Type
1	A	310	ALA

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	378/405 (93%)	359 (95%)	19 (5%)	22	15

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	75	THR
1	A	79	LYS
1	A	116	LYS
1	A	119[A]	ASP
1	A	119[B]	ASP
1	A	120	LYS
1	A	135	ASN
1	A	155	ASP
1	A	164	THR
1	A	209	ASP
1	A	211	THR
1	A	265	ARG
1	A	306	THR
1	A	330	LEU
1	A	338	VAL
1	A	346	PRO
1	A	351	ARG
1	A	437	LEU
1	A	466	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	HIS

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Mol	Chain	Res	Type
1	A	55	HIS
1	A	65	GLN
1	A	66	GLN
1	A	68	GLN
1	A	91	GLN
1	A	136	ASN
1	A	294	HIS
1	A	348	HIS
1	A	371	HIS
1	A	391	ASN
1	A	416	ASN
1	A	438	HIS
1	A	443	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](#)

Of 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 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	UTP	A	502	-	29,30,30	1.50	5 (17%)	43,47,47	1.69	7 (16%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	UTP	A	501	2	29,30,30	1.65	4 (13%)	43,47,47	1.88	6 (13%)
3	UTP	A	503	-	29,30,30	1.63	6 (20%)	43,47,47	1.54	5 (11%)

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	UTP	A	502	-	-	0/22/38/38	0/2/2/2
3	UTP	A	501	2	-	4/22/38/38	0/2/2/2
3	UTP	A	503	-	-	10/22/38/38	0/2/2/2

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	503	UTP	O4-C4	5.98	1.36	1.24
3	A	501	UTP	O4-C4	5.82	1.35	1.24
3	A	502	UTP	O4-C4	5.21	1.34	1.24
3	A	501	UTP	C2-N1	-2.86	1.34	1.38
3	A	503	UTP	PB-O3B	2.82	1.62	1.59
3	A	501	UTP	PA-O3A	2.79	1.62	1.59
3	A	501	UTP	C6-C5	2.44	1.40	1.35
3	A	502	UTP	PB-O3A	2.35	1.62	1.59
3	A	503	UTP	PA-O3A	2.32	1.62	1.59
3	A	503	UTP	C2-N1	-2.31	1.34	1.38
3	A	503	UTP	PB-O3A	2.24	1.61	1.59
3	A	502	UTP	C5-C4	-2.18	1.39	1.43
3	A	502	UTP	C6-C5	2.16	1.40	1.35
3	A	502	UTP	PA-O3A	2.16	1.61	1.59
3	A	503	UTP	C6-C5	2.02	1.39	1.35

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	UTP	C4-N3-C2	-5.89	119.29	126.61
3	A	501	UTP	N3-C2-N1	5.61	122.19	114.89
3	A	503	UTP	C4-N3-C2	-5.16	120.21	126.61
3	A	502	UTP	C4-N3-C2	-4.89	120.54	126.61
3	A	502	UTP	N3-C2-N1	4.38	120.60	114.89
3	A	502	UTP	C5-C4-N3	4.23	120.72	114.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	503	UTP	C5-C4-N3	4.21	120.69	114.80
3	A	501	UTP	C5-C4-N3	4.12	120.57	114.80
3	A	501	UTP	O2-C2-N1	-3.99	117.60	122.80
3	A	503	UTP	N3-C2-N1	3.97	120.06	114.89
3	A	502	UTP	O4-C4-C5	-3.63	118.89	125.16
3	A	501	UTP	O4-C4-C5	-2.87	120.21	125.16
3	A	503	UTP	O4-C4-C5	-2.81	120.31	125.16
3	A	503	UTP	O2-C2-N1	-2.56	119.47	122.80
3	A	502	UTP	C6-C5-C4	-2.23	116.68	119.53
3	A	502	UTP	O2-C2-N1	-2.22	119.90	122.80
3	A	501	UTP	O3A-PB-O2B	-2.18	104.14	110.70
3	A	502	UTP	O2A-PA-O1A	2.03	121.88	112.44

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	501	UTP	PB-O3B-PG-O3G
3	A	503	UTP	C5'-O5'-PA-O2A
3	A	503	UTP	C5'-O5'-PA-O3A
3	A	501	UTP	PA-O3A-PB-O2B
3	A	503	UTP	PG-O3B-PB-O1B
3	A	503	UTP	C5'-O5'-PA-O1A
3	A	501	UTP	PA-O3A-PB-O1B
3	A	503	UTP	PG-O3B-PB-O2B
3	A	503	UTP	PB-O3B-PG-O2G
3	A	501	UTP	PB-O3B-PG-O2G
3	A	503	UTP	PB-O3B-PG-O1G
3	A	503	UTP	PB-O3B-PG-O3G
3	A	503	UTP	PA-O3A-PB-O2B
3	A	503	UTP	PA-O3A-PB-O1B

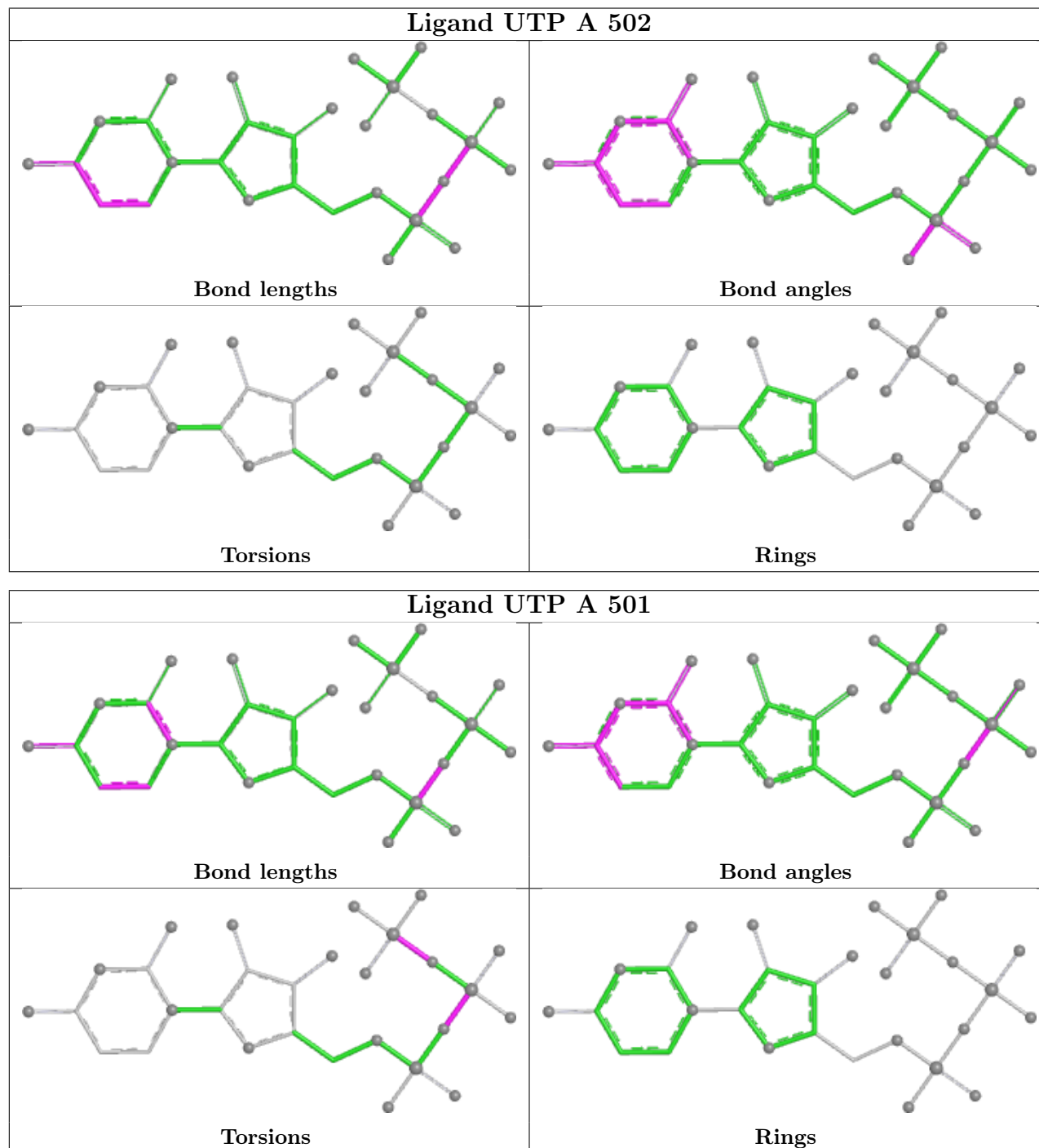
There are no ring outliers.

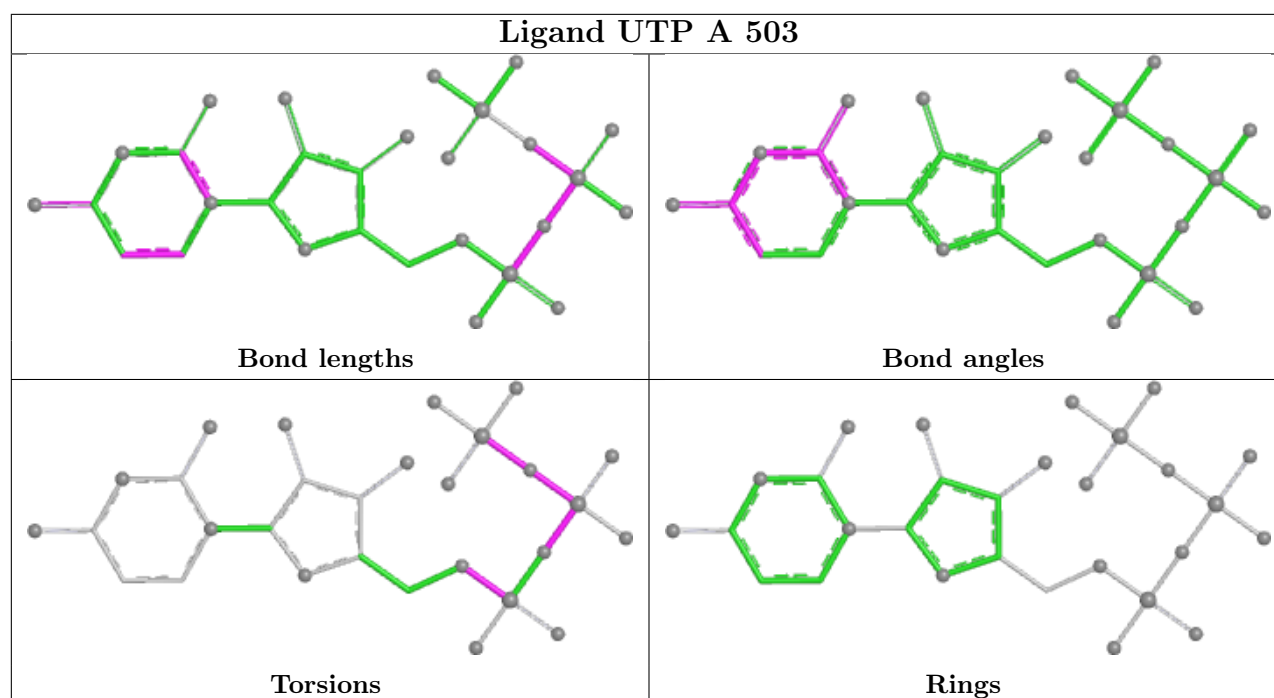
2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502	UTP	3	0
3	A	503	UTP	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.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	444/468 (94%)	0.59	40 (9%) 15 15	21, 31, 47, 59	4 (0%)

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	213	TYR	8.4
1	A	212	THR	7.0
1	A	312	SER	5.5
1	A	310	ALA	5.4
1	A	311	GLY	4.7
1	A	163	TYR	4.6
1	A	30	ASN	4.4
1	A	308	VAL	3.9
1	A	153	ALA	3.8
1	A	211	THR	3.8
1	A	199	GLY	3.7
1	A	313	GLY	3.6
1	A	309	GLY	3.3
1	A	152	PHE	3.2
1	A	112	GLU	3.2
1	A	307	ASN	3.0
1	A	214	GLY	3.0
1	A	314	ALA	3.0
1	A	34	ASP	2.8
1	A	164	THR	2.8
1	A	154	ASN	2.8
1	A	161	ALA	2.8
1	A	200	ARG	2.7
1	A	31	PRO	2.7
1	A	466	GLU	2.7
1	A	306	THR	2.6
1	A	347	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	345	HIS	2.5
1	A	151	ARG	2.5
1	A	210	ARG	2.5
1	A	142	GLU	2.3
1	A	305	ALA	2.3
1	A	160	ARG	2.3
1	A	458	ARG	2.1
1	A	182	ARG	2.1
1	A	216	ARG	2.1
1	A	409	LYS	2.1
1	A	315	MET	2.1
1	A	135	ASN	2.0
1	A	32	SER	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](#)

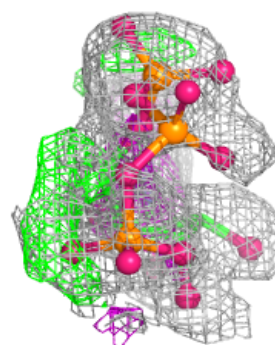
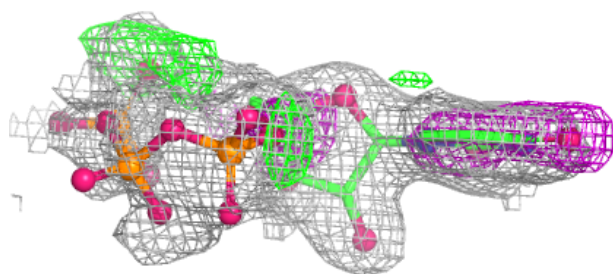
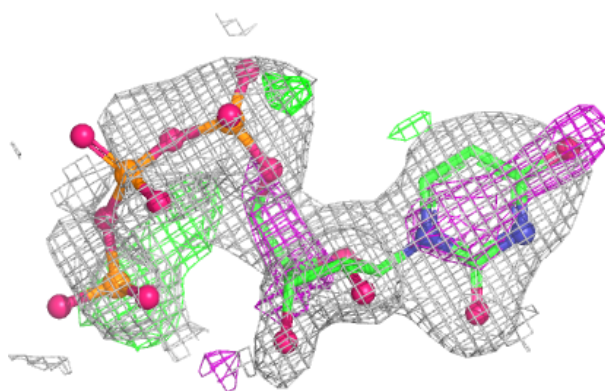
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	UTP	A	503	29/29	0.71	0.23	43,48,57,58	12
3	UTP	A	502	29/29	0.91	0.12	44,47,50,51	4
3	UTP	A	501	29/29	0.98	0.06	26,29,32,33	0
2	MN	A	488	1/1	0.99	0.15	21,21,21,21	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.

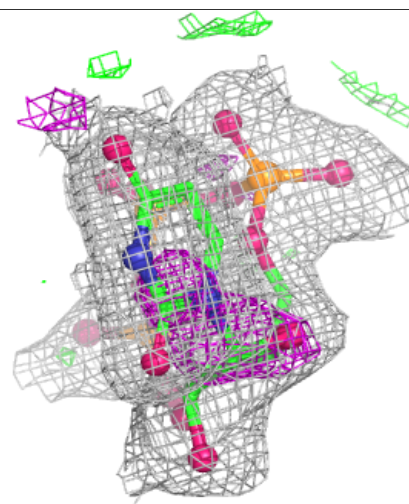
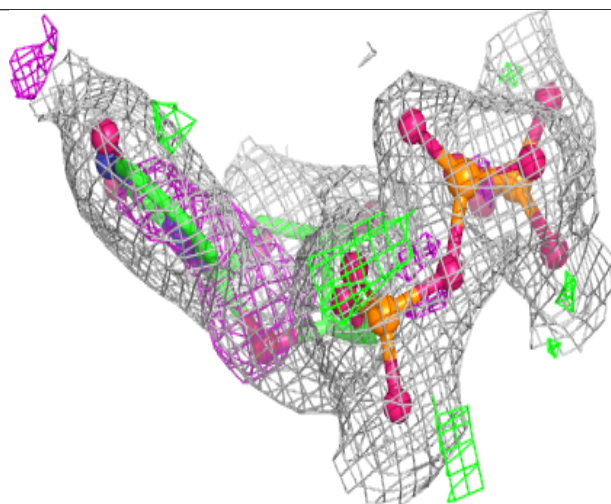
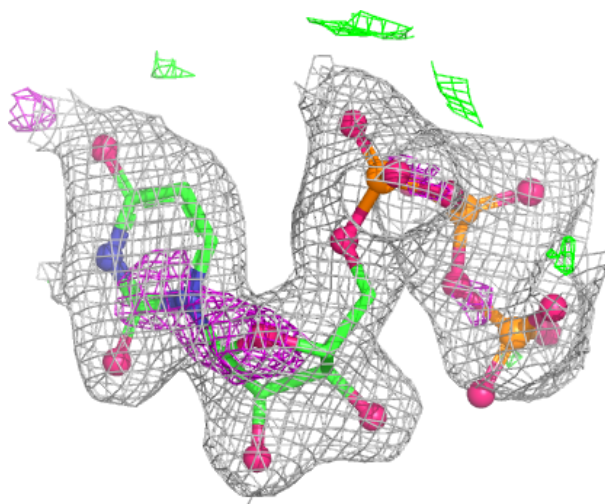
Electron density around UTP A 503:

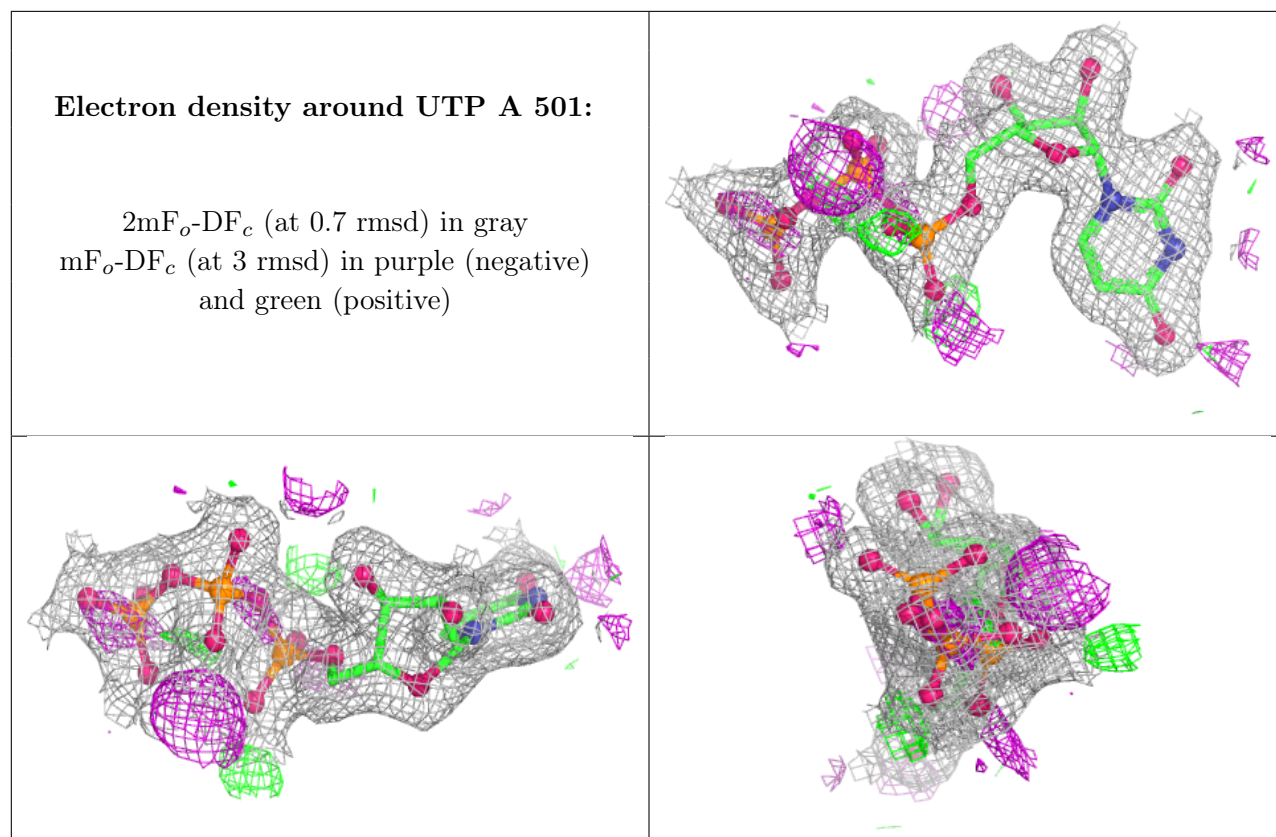
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around UTP 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](#)

There are no such residues in this entry.