



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:05 PM UTC

PDB ID : 5Z09 / pdb_00005z09
Title : ST0452(Y97N)-UTP binding form
Authors : Honda, Y.; Nakano, S.; Ito, S.; Dadashipour, M.; Zhang, Z.; Kawarabayasi, Y.
Deposited on : 2017-12-19
Resolution : 2.91 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

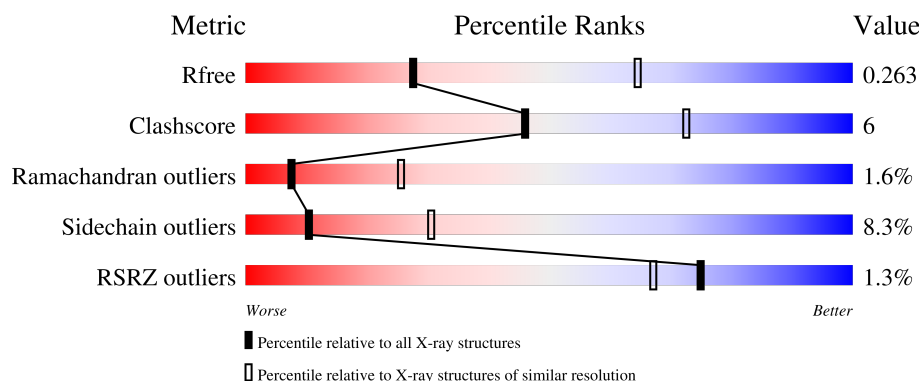
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.91 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	2995 (2.94-2.90)
Clashscore	190562	3213 (2.94-2.90)
Ramachandran outliers	187476	3128 (2.94-2.90)
Sidechain outliers	187428	3130 (2.94-2.90)
RSRZ outliers	180081	2995 (2.94-2.90)

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	409	% 80% 15% ..
1	B	409	81% 14% ...
1	C	409	% 82% 14% ..
1	D	409	% 82% 14% ..
1	E	409	% 74% 13% • 10%

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Mol	Chain	Length	Quality of chain
1	F	409	2% 71% 13% • • 11%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 18308 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 Dual sugar-1-phosphate nucleotidyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	401	Total	C	N	O	S	0	0	0
			3130	2013	522	589	6			
1	B	401	Total	C	N	O	S	0	0	0
			3118	2006	521	585	6			
1	C	401	Total	C	N	O	S	0	0	0
			3122	2008	521	587	6			
1	D	401	Total	C	N	O	S	0	0	0
			3122	2008	521	587	6			
1	E	369	Total	C	N	O	S	0	0	0
			2873	1847	483	538	5			
1	F	364	Total	C	N	O	S	0	0	0
			2827	1820	474	528	5			

There are 54 discrepancies between the modelled and reference sequences:

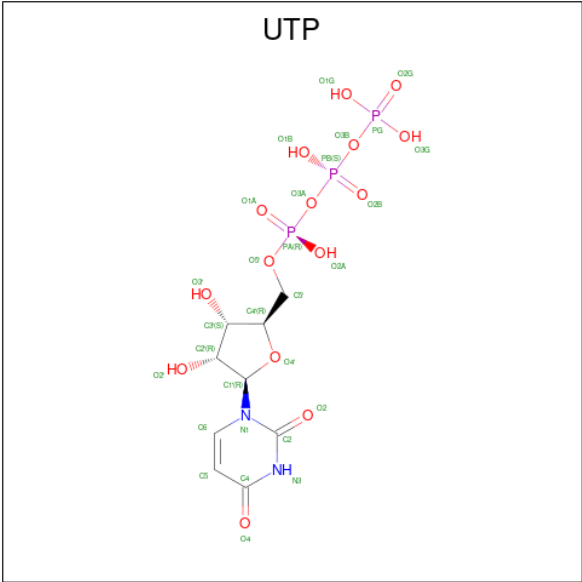
Chain	Residue	Modelled	Actual	Comment	Reference
A	97	ASN	TYR	engineered mutation	UNP Q975F9
A	402	LEU	-	expression tag	UNP Q975F9
A	403	GLU	-	expression tag	UNP Q975F9
A	404	HIS	-	expression tag	UNP Q975F9
A	405	HIS	-	expression tag	UNP Q975F9
A	406	HIS	-	expression tag	UNP Q975F9
A	407	HIS	-	expression tag	UNP Q975F9
A	408	HIS	-	expression tag	UNP Q975F9
A	409	HIS	-	expression tag	UNP Q975F9
B	97	ASN	TYR	engineered mutation	UNP Q975F9
B	402	LEU	-	expression tag	UNP Q975F9
B	403	GLU	-	expression tag	UNP Q975F9
B	404	HIS	-	expression tag	UNP Q975F9
B	405	HIS	-	expression tag	UNP Q975F9
B	406	HIS	-	expression tag	UNP Q975F9
B	407	HIS	-	expression tag	UNP Q975F9
B	408	HIS	-	expression tag	UNP Q975F9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	409	HIS	-	expression tag	UNP Q975F9
C	97	ASN	TYR	engineered mutation	UNP Q975F9
C	402	LEU	-	expression tag	UNP Q975F9
C	403	GLU	-	expression tag	UNP Q975F9
C	404	HIS	-	expression tag	UNP Q975F9
C	405	HIS	-	expression tag	UNP Q975F9
C	406	HIS	-	expression tag	UNP Q975F9
C	407	HIS	-	expression tag	UNP Q975F9
C	408	HIS	-	expression tag	UNP Q975F9
C	409	HIS	-	expression tag	UNP Q975F9
D	97	ASN	TYR	engineered mutation	UNP Q975F9
D	402	LEU	-	expression tag	UNP Q975F9
D	403	GLU	-	expression tag	UNP Q975F9
D	404	HIS	-	expression tag	UNP Q975F9
D	405	HIS	-	expression tag	UNP Q975F9
D	406	HIS	-	expression tag	UNP Q975F9
D	407	HIS	-	expression tag	UNP Q975F9
D	408	HIS	-	expression tag	UNP Q975F9
D	409	HIS	-	expression tag	UNP Q975F9
E	97	ASN	TYR	engineered mutation	UNP Q975F9
E	402	LEU	-	expression tag	UNP Q975F9
E	403	GLU	-	expression tag	UNP Q975F9
E	404	HIS	-	expression tag	UNP Q975F9
E	405	HIS	-	expression tag	UNP Q975F9
E	406	HIS	-	expression tag	UNP Q975F9
E	407	HIS	-	expression tag	UNP Q975F9
E	408	HIS	-	expression tag	UNP Q975F9
E	409	HIS	-	expression tag	UNP Q975F9
F	97	ASN	TYR	engineered mutation	UNP Q975F9
F	402	LEU	-	expression tag	UNP Q975F9
F	403	GLU	-	expression tag	UNP Q975F9
F	404	HIS	-	expression tag	UNP Q975F9
F	405	HIS	-	expression tag	UNP Q975F9
F	406	HIS	-	expression tag	UNP Q975F9
F	407	HIS	-	expression tag	UNP Q975F9
F	408	HIS	-	expression tag	UNP Q975F9
F	409	HIS	-	expression tag	UNP Q975F9

- Molecule 2 is URIDINE 5'-TRIPHOSPHATE (CCD ID: UTP) (formula: $C_9H_{15}N_2O_{15}P_3$).

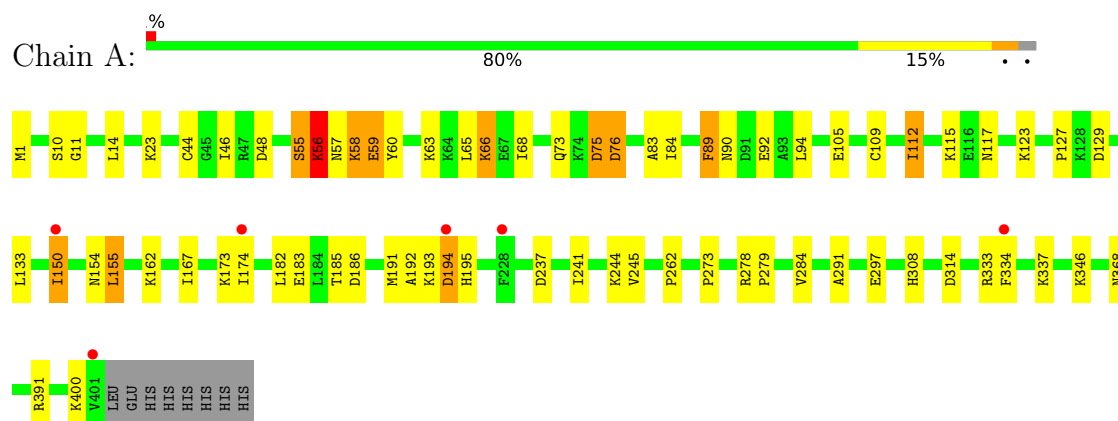


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			29	9	2	15	3		
2	B	1	Total	C	N	O	P	0	0
			29	9	2	15	3		
2	C	1	Total	C	N	O	P	0	0
			29	9	2	15	3		
2	D	1	Total	C	N	O	P	0	0
			29	9	2	15	3		

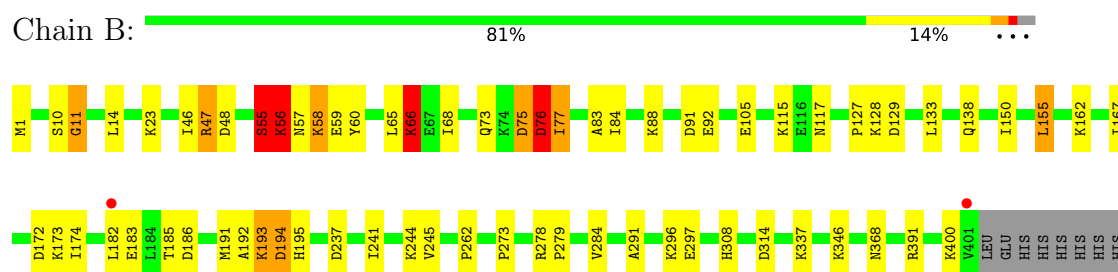
3 Residue-property plots [i](#)

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.

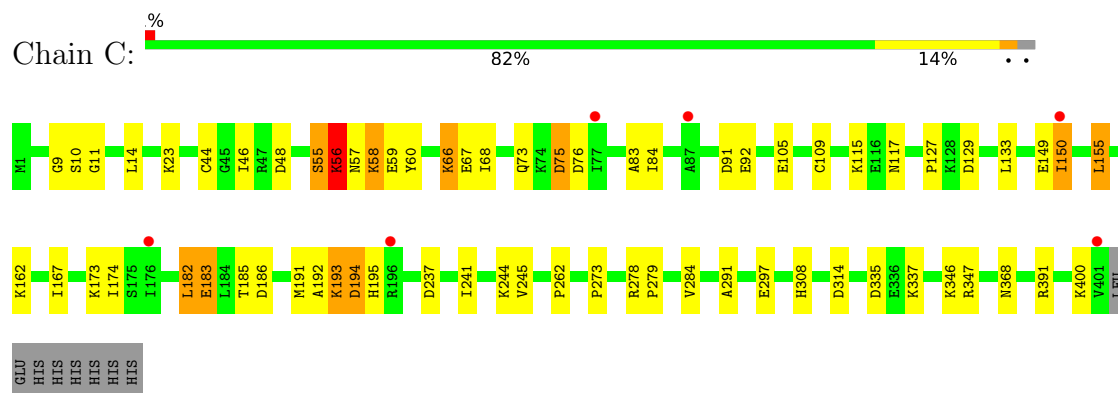
- Molecule 1: Dual sugar-1-phosphate nucleotidyltransferase



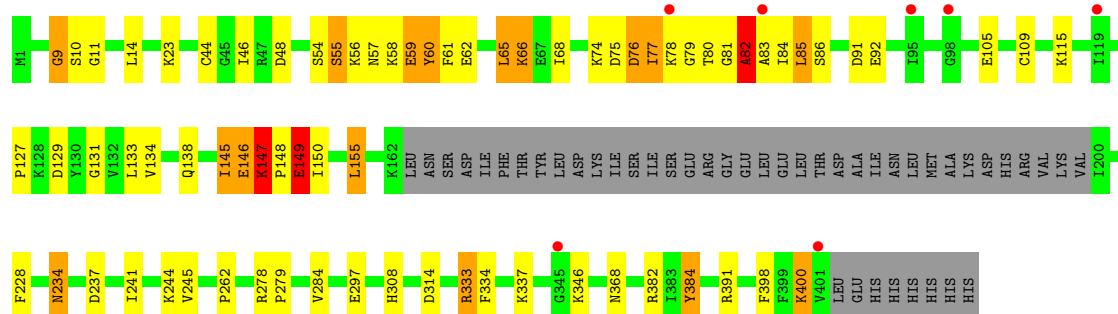
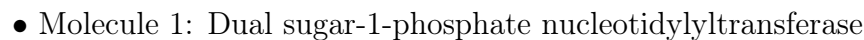
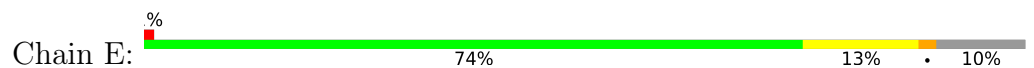
- Molecule 1: Dual sugar-1-phosphate nucleotidyltransferase



- Molecule 1: Dual sugar-1-phosphate nucleotidyltransferase



Chain D: 82% 14% ..



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	210.66Å 88.42Å 179.34Å 90.00° 121.59° 90.00°	Depositor
Resolution (Å)	152.76 – 2.91 152.76 – 2.91	Depositor EDS
% Data completeness (in resolution range)	98.5 (152.76-2.91) 98.6 (152.76-2.91)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.209 , 0.267 0.210 , 0.263	Depositor DCC
R_{free} test set	3064 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	69.6	Xtriage
Anisotropy	0.084	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 56.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	18308	wwPDB-VP
Average B, all atoms (Å ²)	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.24% 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: 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	0.53	0/3183	0.73	0/4297
1	B	0.51	0/3171	0.70	0/4283
1	C	0.48	0/3175	0.68	0/4288
1	D	0.47	0/3175	0.70	1/4288 (0.0%)
1	E	0.49	0/2920	0.76	5/3941 (0.1%)
1	F	0.47	0/2876	0.73	3/3882 (0.1%)
All	All	0.49	0/18500	0.72	9/24979 (0.0%)

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	9
1	B	0	6
1	C	0	6
1	D	0	7
1	E	0	8
1	F	0	4
All	All	0	40

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	8	ALA	N-CA-C	-7.92	97.69	109.81
1	F	384	TYR	CB-CA-C	6.36	119.40	109.42
1	E	84	ILE	CA-C-N	6.31	133.06	121.70
1	E	84	ILE	C-N-CA	6.31	133.06	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	83	ALA	N-CA-C	5.85	118.13	108.48
1	D	195	HIS	N-CA-C	5.68	120.24	112.68
1	E	334	PHE	N-CA-C	-5.39	101.87	109.96
1	F	147	LYS	CA-C-N	-5.36	108.61	120.56
1	F	147	LYS	C-N-CA	-5.36	108.61	120.56

There are no chirality outliers.

All (40) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	194	ASP	Peptide
1	A	55	SER	Peptide
1	A	56	LYS	Peptide
1	A	58	LYS	Peptide
1	A	59	GLU	Peptide,Mainchain
1	A	65	LEU	Peptide,Mainchain
1	A	89	PHE	Peptide
1	B	194	ASP	Peptide
1	B	55	SER	Peptide
1	B	56	LYS	Peptide
1	B	58	LYS	Peptide
1	B	59	GLU	Peptide
1	B	66	LYS	Peptide
1	C	194	ASP	Peptide
1	C	55	SER	Peptide
1	C	56	LYS	Peptide
1	C	58	LYS	Peptide
1	C	59	GLU	Peptide
1	C	9	GLY	Peptide
1	D	195	HIS	Peptide
1	D	196	ARG	Peptide
1	D	55	SER	Peptide
1	D	56	LYS	Peptide
1	D	58	LYS	Peptide
1	D	59	GLU	Peptide
1	D	65	LEU	Peptide
1	E	10	SER	Peptide
1	E	55	SER	Peptide
1	E	56	LYS	Peptide
1	E	58	LYS	Peptide
1	E	59	GLU	Peptide
1	E	65	LEU	Peptide

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Mol	Chain	Res	Type	Group
1	E	74	LYS	Peptide
1	E	9	GLY	Peptide
1	F	65	LEU	Peptide
1	F	81	GLY	Peptide
1	F	82	ALA	Peptide
1	F	9	GLY	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	3130	0	3230	39	0
1	B	3118	0	3211	31	0
1	C	3122	0	3215	28	0
1	D	3122	0	3215	23	0
1	E	2873	0	2943	47	0
1	F	2827	0	2911	58	0
2	A	29	0	11	0	0
2	B	29	0	11	1	0
2	C	29	0	11	1	0
2	D	29	0	11	0	0
All	All	18308	0	18769	216	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 6.

All (216) 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:123:LYS:NZ	1:E:69:SER:OG	1.65	1.28
1:E:65:LEU:HA	1:E:66:LYS:HB2	1.23	1.19
1:D:65:LEU:HA	1:D:66:LYS:HB2	1.22	1.15
1:E:83:ALA:HB3	1:E:84:ILE:HA	1.28	1.12
1:F:65:LEU:HA	1:F:66:LYS:HB2	1.23	1.09
1:A:1:MET:HE3	1:A:112:ILE:HG13	1.33	1.06
1:F:59:GLU:HA	1:F:61:PHE:N	1.71	1.04

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:65:LEU:HA	1:E:66:LYS:CB	1.97	0.95
1:D:65:LEU:HA	1:D:66:LYS:CB	1.97	0.93
1:F:65:LEU:HA	1:F:66:LYS:CB	1.97	0.93
1:F:59:GLU:CB	1:F:62:GLU:H	1.83	0.92
1:B:75:ASP:O	1:B:77:ILE:N	2.04	0.89
1:F:58:LYS:O	1:F:59:GLU:CB	2.20	0.88
1:F:59:GLU:HA	1:F:60:TYR:C	1.97	0.86
1:F:134:VAL:CG2	1:F:145:ILE:HD11	2.06	0.85
1:F:148:PRO:O	1:F:150:ILE:N	2.08	0.85
1:F:146:GLU:O	1:F:147:LYS:HG3	1.78	0.83
1:F:82:ALA:HB1	1:F:83:ALA:C	2.05	0.82
1:F:59:GLU:CB	1:F:62:GLU:CG	2.58	0.81
1:E:199:VAL:HG13	1:E:199:VAL:O	1.78	0.80
1:C:182:LEU:HD12	1:C:182:LEU:O	1.83	0.78
1:A:1:MET:HE3	1:A:112:ILE:CG1	2.14	0.76
1:E:85:LEU:HD23	1:E:85:LEU:C	2.11	0.75
1:E:106:LYS:HD2	1:F:228:PHE:CD2	2.21	0.75
1:F:54:SER:O	1:F:57:ASN:HB2	1.87	0.74
1:E:83:ALA:HB3	1:E:84:ILE:CA	2.15	0.74
1:F:59:GLU:CB	1:F:62:GLU:HG2	2.19	0.72
1:A:123:LYS:NZ	1:E:69:SER:HG	1.86	0.71
1:E:83:ALA:CB	1:E:84:ILE:HA	2.10	0.71
1:F:134:VAL:HG21	1:F:145:ILE:HD11	1.73	0.69
1:E:83:ALA:H	1:E:84:ILE:HG13	1.57	0.69
1:B:193:LYS:N	1:B:194:ASP:HA	2.08	0.69
1:A:1:MET:HG2	1:A:112:ILE:HG12	1.75	0.68
1:A:193:LYS:N	1:A:194:ASP:HA	2.09	0.68
1:E:84:ILE:HG23	1:E:85:LEU:HB3	1.73	0.68
1:E:199:VAL:O	1:E:199:VAL:CG1	2.41	0.67
1:C:193:LYS:N	1:C:194:ASP:HA	2.09	0.66
1:F:9:GLY:O	1:F:10:SER:OG	2.12	0.66
1:F:333:ARG:HD3	1:F:334:PHE:H	1.62	0.65
1:F:59:GLU:HA	1:F:61:PHE:H	1.61	0.65
1:F:83:ALA:HA	1:F:85:LEU:H	1.61	0.65
1:F:85:LEU:C	1:F:85:LEU:HD22	2.22	0.65
1:A:150:ILE:CG2	1:F:234:ASN:HB3	2.27	0.65
1:F:148:PRO:O	1:F:149:GLU:C	2.39	0.64
1:B:75:ASP:C	1:B:77:ILE:H	2.03	0.64
1:F:150:ILE:HD12	1:F:150:ILE:O	1.97	0.64
1:E:334:PHE:O	1:E:335:ASP:CG	2.42	0.62
1:E:84:ILE:HA	1:E:85:LEU:HB3	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:65:LEU:CA	1:F:66:LYS:CB	2.76	0.61
1:E:83:ALA:HB3	1:E:85:LEU:HB2	1.81	0.60
1:F:148:PRO:C	1:F:150:ILE:N	2.59	0.60
1:A:1:MET:HG2	1:A:112:ILE:CG1	2.30	0.60
1:B:117:ASN:ND2	1:B:191:MET:SD	2.76	0.59
1:E:52:ILE:HD13	1:E:71:VAL:CG1	2.34	0.58
1:E:84:ILE:HA	1:E:85:LEU:CB	2.33	0.58
1:E:81:GLY:N	1:E:82:ALA:HB2	2.19	0.58
1:E:85:LEU:HD23	1:E:85:LEU:O	2.04	0.57
1:E:65:LEU:CA	1:E:66:LYS:CB	2.76	0.56
1:A:58:LYS:HB3	1:A:59:GLU:HG2	1.88	0.56
1:E:65:LEU:CA	1:E:66:LYS:HB2	2.14	0.56
1:F:85:LEU:HD13	1:F:86:SER:N	2.21	0.55
1:C:182:LEU:HD12	1:C:182:LEU:C	2.30	0.55
1:E:6:LEU:HD13	1:E:83:ALA:HB2	1.87	0.55
1:D:193:LYS:C	1:D:194:ASP:CG	2.74	0.55
1:D:65:LEU:CA	1:D:66:LYS:CB	2.75	0.55
1:D:174:ILE:HG12	1:D:186:ASP:HB3	1.89	0.55
1:C:117:ASN:ND2	1:C:191:MET:SD	2.81	0.54
1:A:59:GLU:O	1:A:63:LYS:HG2	2.08	0.54
1:A:117:ASN:ND2	1:A:191:MET:SD	2.81	0.54
1:C:241:ILE:HG23	1:C:245:VAL:HG21	1.90	0.54
1:E:84:ILE:CA	1:E:85:LEU:HB3	2.38	0.53
1:F:146:GLU:HG3	1:F:147:LYS:HG2	1.89	0.53
1:F:59:GLU:CB	1:F:62:GLU:HG3	2.37	0.53
1:B:278:ARG:HB3	1:B:279:PRO:HD2	1.91	0.53
1:B:241:ILE:HG23	1:B:245:VAL:HG21	1.90	0.53
1:E:84:ILE:CG2	1:E:85:LEU:HB3	2.39	0.53
1:E:278:ARG:HB3	1:E:279:PRO:HD2	1.91	0.53
1:D:117:ASN:ND2	1:D:191:MET:SD	2.82	0.53
1:E:241:ILE:HG23	1:E:245:VAL:HG21	1.90	0.53
1:A:278:ARG:HB3	1:A:279:PRO:HD2	1.91	0.52
1:C:192:ALA:O	1:C:194:ASP:HB3	2.09	0.52
1:D:241:ILE:HG23	1:D:245:VAL:HG21	1.90	0.52
1:F:241:ILE:HG23	1:F:245:VAL:HG21	1.91	0.52
1:B:174:ILE:HG12	1:B:186:ASP:HB3	1.92	0.52
1:A:1:MET:CE	1:A:112:ILE:HG13	2.24	0.52
1:D:195:HIS:CB	1:D:196:ARG:HA	2.40	0.52
1:A:192:ALA:O	1:A:194:ASP:HB3	2.09	0.52
1:A:14:LEU:HD22	1:A:23:LYS:HD3	1.93	0.51
1:A:241:ILE:HG23	1:A:245:VAL:HG21	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:174:ILE:HG12	1:A:186:ASP:HB3	1.91	0.51
1:C:278:ARG:HB3	1:C:279:PRO:HD2	1.93	0.51
1:F:134:VAL:HG23	1:F:145:ILE:HD11	1.92	0.51
1:F:278:ARG:HB3	1:F:279:PRO:HD2	1.92	0.51
1:F:55:SER:OG	1:F:58:LYS:HB2	2.11	0.50
1:B:192:ALA:O	1:B:194:ASP:HB3	2.11	0.50
1:F:382:ARG:HD2	1:F:398:PHE:CE2	2.46	0.50
1:C:174:ILE:HG12	1:C:186:ASP:HB3	1.92	0.50
1:D:195:HIS:HB2	1:D:196:ARG:HA	1.94	0.50
1:E:85:LEU:C	1:E:85:LEU:CD2	2.80	0.50
1:A:56:LYS:HG2	1:A:57:ASN:HA	1.94	0.50
1:F:146:GLU:O	1:F:147:LYS:CG	2.55	0.50
1:D:14:LEU:HD22	1:D:23:LYS:HD3	1.94	0.49
1:A:1:MET:HG2	1:A:112:ILE:CD1	2.42	0.49
1:C:14:LEU:HD22	1:C:23:LYS:HD3	1.94	0.49
1:A:150:ILE:HG22	1:F:234:ASN:HB3	1.92	0.49
1:B:56:LYS:HG3	1:C:150:ILE:HG12	1.93	0.49
1:D:278:ARG:HB3	1:D:279:PRO:HD2	1.93	0.49
1:E:82:ALA:CB	1:E:84:ILE:HD11	2.43	0.49
2:C:501:UTP:H3'	2:C:501:UTP:O2A	2.12	0.49
1:F:54:SER:O	1:F:57:ASN:CB	2.58	0.49
1:B:56:LYS:HG2	1:B:57:ASN:HA	1.95	0.49
1:D:244:LYS:O	1:D:262:PRO:HA	2.12	0.49
1:E:14:LEU:HD22	1:E:23:LYS:HD3	1.95	0.48
1:A:66:LYS:H	1:A:66:LYS:HD2	1.78	0.48
1:B:14:LEU:HD22	1:B:23:LYS:HD3	1.95	0.48
1:D:56:LYS:HG2	1:D:57:ASN:HA	1.95	0.48
1:A:297:GLU:O	1:A:314:ASP:HA	2.13	0.48
1:F:59:GLU:CB	1:F:62:GLU:CB	2.92	0.48
1:A:1:MET:HG2	1:A:112:ILE:HD11	1.95	0.48
1:C:56:LYS:HG2	1:C:57:ASN:HA	1.95	0.48
1:F:65:LEU:CA	1:F:66:LYS:HB2	2.14	0.48
1:C:244:LYS:O	1:C:262:PRO:HA	2.14	0.48
1:A:150:ILE:CG2	1:F:234:ASN:CB	2.92	0.48
1:B:244:LYS:O	1:B:262:PRO:HA	2.14	0.48
1:F:76:ASP:O	1:F:77:ILE:HB	2.14	0.48
1:B:193:LYS:N	1:B:194:ASP:CA	2.77	0.47
1:F:14:LEU:HD22	1:F:23:LYS:HD3	1.95	0.47
1:F:148:PRO:C	1:F:150:ILE:H	2.23	0.47
1:A:244:LYS:O	1:A:262:PRO:HA	2.14	0.47
1:B:278:ARG:HB3	1:B:279:PRO:CD	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:297:GLU:O	1:B:314:ASP:HA	2.15	0.47
1:F:244:LYS:O	1:F:262:PRO:HA	2.14	0.47
1:F:82:ALA:HA	1:F:83:ALA:HB3	1.97	0.47
1:E:56:LYS:HG2	1:E:57:ASN:HA	1.96	0.46
1:F:146:GLU:OE2	1:F:147:LYS:N	2.48	0.46
1:A:46:ILE:O	1:A:48:ASP:N	2.48	0.46
1:C:193:LYS:N	1:C:194:ASP:CA	2.79	0.46
1:E:83:ALA:N	1:E:84:ILE:HG13	2.26	0.46
1:D:297:GLU:O	1:D:314:ASP:HA	2.14	0.46
1:C:297:GLU:O	1:C:314:ASP:HA	2.16	0.45
1:C:278:ARG:HB3	1:C:279:PRO:CD	2.46	0.45
1:D:333:ARG:NH1	1:D:335:ASP:OD2	2.49	0.45
1:E:278:ARG:HB3	1:E:279:PRO:CD	2.46	0.45
1:A:193:LYS:N	1:A:194:ASP:CA	2.78	0.45
1:F:297:GLU:O	1:F:314:ASP:HA	2.16	0.45
1:B:76:ASP:C	1:B:77:ILE:HD13	2.42	0.45
1:C:46:ILE:O	1:C:48:ASP:N	2.49	0.45
1:E:82:ALA:HB3	1:E:84:ILE:HD11	1.98	0.45
1:E:244:LYS:O	1:E:262:PRO:HA	2.16	0.45
1:F:79:GLY:HA3	1:F:80:THR:HA	1.70	0.45
1:A:1:MET:CE	1:A:112:ILE:CG1	2.92	0.45
1:F:54:SER:O	1:F:57:ASN:N	2.50	0.45
1:A:278:ARG:HB3	1:A:279:PRO:CD	2.46	0.45
1:E:46:ILE:O	1:E:48:ASP:N	2.48	0.45
1:E:297:GLU:O	1:E:314:ASP:HA	2.17	0.45
1:D:278:ARG:HB3	1:D:279:PRO:CD	2.47	0.44
1:F:278:ARG:HB3	1:F:279:PRO:CD	2.46	0.44
1:B:1:MET:O	1:B:47:ARG:NH2	2.50	0.44
1:B:65:LEU:C	1:B:66:LYS:HE2	2.43	0.44
1:B:55:SER:HB2	1:C:149:GLU:OE1	2.18	0.44
1:F:127:PRO:O	1:F:129:ASP:O	2.36	0.44
1:E:127:PRO:O	1:E:129:ASP:O	2.36	0.44
1:A:194:ASP:HB2	1:A:195:HIS:O	2.18	0.43
1:A:154:ASN:ND2	1:E:48:ASP:HB2	2.34	0.43
1:B:66:LYS:HE2	1:B:66:LYS:N	2.33	0.43
1:F:82:ALA:HB1	1:F:83:ALA:CA	2.48	0.43
1:D:127:PRO:O	1:D:129:ASP:O	2.37	0.43
1:A:127:PRO:O	1:A:129:ASP:O	2.36	0.43
1:A:133:LEU:HD12	1:A:155:LEU:HD13	2.00	0.43
1:B:46:ILE:O	1:B:48:ASP:N	2.49	0.43
1:C:127:PRO:O	1:C:129:ASP:O	2.36	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:74:LYS:HB2	1:F:75:ASP:HB2	2.01	0.43
1:B:73:GLN:HG3	1:B:83:ALA:HA	2.01	0.43
1:B:127:PRO:O	1:B:129:ASP:O	2.37	0.43
1:D:46:ILE:O	1:D:48:ASP:N	2.49	0.43
1:D:333:ARG:HD2	1:D:339:VAL:HA	2.00	0.43
1:B:56:LYS:HA	1:B:58:LYS:H	1.83	0.42
1:E:52:ILE:CD1	1:E:87:ALA:CB	2.97	0.42
1:F:131:GLY:HA2	1:F:146:GLU:HA	2.01	0.42
1:F:384:TYR:HB2	1:F:400:LYS:HA	2.00	0.42
1:B:11:GLY:HA2	2:B:501:UTP:O3G	2.19	0.42
1:D:296:LYS:O	1:D:297:GLU:C	2.62	0.42
1:E:83:ALA:HB3	1:E:85:LEU:CB	2.48	0.42
1:F:46:ILE:O	1:F:48:ASP:N	2.49	0.42
1:F:133:LEU:HD12	1:F:155:LEU:HD13	2.01	0.42
1:C:73:GLN:HG3	1:C:83:ALA:HA	2.01	0.42
1:C:133:LEU:HD12	1:C:155:LEU:HD13	2.01	0.42
1:E:52:ILE:HD11	1:E:87:ALA:CB	2.49	0.42
1:B:88:LYS:HZ3	1:C:347:ARG:HB3	1.85	0.42
1:B:273:PRO:O	1:B:291:ALA:HA	2.20	0.42
1:D:133:LEU:HD12	1:D:155:LEU:HD13	2.02	0.42
1:E:133:LEU:HD12	1:E:155:LEU:HD13	2.01	0.41
1:E:83:ALA:CB	1:E:84:ILE:CA	2.80	0.41
1:A:44:CYS:SG	1:A:109:CYS:N	2.93	0.41
1:B:133:LEU:HD12	1:B:155:LEU:HD13	2.01	0.41
1:A:73:GLN:HG3	1:A:83:ALA:HA	2.02	0.41
1:A:94:LEU:HD13	1:A:112:ILE:HD12	2.01	0.41
1:A:89:PHE:HB2	1:A:90:ASN:HB2	2.03	0.41
1:A:273:PRO:O	1:A:291:ALA:HA	2.21	0.41
1:B:296:LYS:O	1:B:297:GLU:C	2.63	0.41
1:C:44:CYS:SG	1:C:109:CYS:N	2.94	0.41
1:C:194:ASP:HB2	1:C:195:HIS:O	2.21	0.41
1:E:54:SER:O	1:E:57:ASN:N	2.54	0.41
1:E:106:LYS:HB2	1:E:106:LYS:HE2	1.82	0.41
1:F:77:ILE:HG23	1:F:78:LYS:N	2.36	0.41
1:E:52:ILE:CD1	1:E:71:VAL:CG1	2.99	0.40
1:F:44:CYS:SG	1:F:109:CYS:N	2.94	0.40
1:D:211:LYS:HD2	1:D:213:TRP:CZ2	2.56	0.40
1:C:66:LYS:HA	1:C:67:GLU:HA	1.87	0.40
1:C:75:ASP:OD1	1:C:75:ASP:N	2.52	0.40
1:D:73:GLN:HG3	1:D:83:ALA:HA	2.02	0.40
1:B:172:ASP:OD2	1:C:335:ASP:HB2	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:56:LYS:HA	1:C:58:LYS:H	1.87	0.40
1:A:75:ASP:OD1	1:A:75:ASP:N	2.52	0.40
1:B:194:ASP:HB2	1:B:195:HIS:O	2.22	0.40
1:C:273:PRO:O	1:C:291:ALA:HA	2.22	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	399/409 (98%)	377 (94%)	19 (5%)	3 (1%)	16	43
1	B	399/409 (98%)	375 (94%)	19 (5%)	5 (1%)	9	30
1	C	399/409 (98%)	377 (94%)	17 (4%)	5 (1%)	9	30
1	D	399/409 (98%)	373 (94%)	20 (5%)	6 (2%)	8	27
1	E	359/409 (88%)	327 (91%)	25 (7%)	7 (2%)	6	22
1	F	360/409 (88%)	326 (91%)	24 (7%)	10 (3%)	4	14
All	All	2315/2454 (94%)	2155 (93%)	124 (5%)	36 (2%)	7	25

All (36) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	76	ASP
1	F	59	GLU
1	F	77	ILE
1	F	82	ALA
1	F	149	GLU
1	A	11	GLY
1	A	60	TYR
1	A	76	ASP

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Mol	Chain	Res	Type
1	B	11	GLY
1	B	60	TYR
1	B	91	ASP
1	C	11	GLY
1	C	91	ASP
1	C	183	GLU
1	D	11	GLY
1	D	66	LYS
1	D	76	ASP
1	D	91	ASP
1	D	194	ASP
1	E	11	GLY
1	E	66	LYS
1	E	85	LEU
1	E	335	ASP
1	F	11	GLY
1	F	66	LYS
1	F	76	ASP
1	F	91	ASP
1	B	75	ASP
1	C	60	TYR
1	D	60	TYR
1	E	60	TYR
1	E	86	SER
1	C	76	ASP
1	F	60	TYR
1	E	75	ASP
1	F	84	ILE

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	343/352 (97%)	313 (91%)	30 (9%)	9	28
1	B	340/352 (97%)	309 (91%)	31 (9%)	9	27
1	C	341/352 (97%)	314 (92%)	27 (8%)	11	33

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	341/352 (97%)	313 (92%)	28 (8%)	10	31
1	E	313/352 (89%)	288 (92%)	25 (8%)	11	32
1	F	307/352 (87%)	284 (92%)	23 (8%)	12	35
All	All	1985/2112 (94%)	1821 (92%)	164 (8%)	10	30

All (164) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	SER
1	A	55	SER
1	A	56	LYS
1	A	66	LYS
1	A	68	ILE
1	A	75	ASP
1	A	76	ASP
1	A	84	ILE
1	A	92	GLU
1	A	105	GLU
1	A	112	ILE
1	A	115	LYS
1	A	150	ILE
1	A	155	LEU
1	A	162	LYS
1	A	167	ILE
1	A	173	LYS
1	A	182	LEU
1	A	183	GLU
1	A	185	THR
1	A	237	ASP
1	A	284	VAL
1	A	308	HIS
1	A	333	ARG
1	A	334	PHE
1	A	337	LYS
1	A	346	LYS
1	A	368	ASN
1	A	391	ARG
1	A	400	LYS
1	B	10	SER
1	B	47	ARG
1	B	55	SER

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Mol	Chain	Res	Type
1	B	56	LYS
1	B	66	LYS
1	B	68	ILE
1	B	76	ASP
1	B	77	ILE
1	B	84	ILE
1	B	92	GLU
1	B	105	GLU
1	B	115	LYS
1	B	128	LYS
1	B	138	GLN
1	B	150	ILE
1	B	155	LEU
1	B	162	LYS
1	B	167	ILE
1	B	173	LYS
1	B	182	LEU
1	B	183	GLU
1	B	185	THR
1	B	193	LYS
1	B	237	ASP
1	B	284	VAL
1	B	308	HIS
1	B	337	LYS
1	B	346	LYS
1	B	368	ASN
1	B	391	ARG
1	B	400	LYS
1	C	10	SER
1	C	55	SER
1	C	56	LYS
1	C	66	LYS
1	C	68	ILE
1	C	75	ASP
1	C	84	ILE
1	C	92	GLU
1	C	105	GLU
1	C	115	LYS
1	C	150	ILE
1	C	155	LEU
1	C	162	LYS
1	C	167	ILE

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Mol	Chain	Res	Type
1	C	173	LYS
1	C	182	LEU
1	C	183	GLU
1	C	185	THR
1	C	193	LYS
1	C	237	ASP
1	C	284	VAL
1	C	308	HIS
1	C	337	LYS
1	C	346	LYS
1	C	368	ASN
1	C	391	ARG
1	C	400	LYS
1	D	10	SER
1	D	55	SER
1	D	56	LYS
1	D	68	ILE
1	D	76	ASP
1	D	84	ILE
1	D	92	GLU
1	D	105	GLU
1	D	115	LYS
1	D	138	GLN
1	D	150	ILE
1	D	155	LEU
1	D	162	LYS
1	D	167	ILE
1	D	173	LYS
1	D	182	LEU
1	D	183	GLU
1	D	185	THR
1	D	197	VAL
1	D	237	ASP
1	D	284	VAL
1	D	308	HIS
1	D	333	ARG
1	D	337	LYS
1	D	346	LYS
1	D	368	ASN
1	D	391	ARG
1	D	400	LYS
1	E	55	SER

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Mol	Chain	Res	Type
1	E	56	LYS
1	E	68	ILE
1	E	74	LYS
1	E	76	ASP
1	E	85	LEU
1	E	88	LYS
1	E	91	ASP
1	E	92	GLU
1	E	105	GLU
1	E	115	LYS
1	E	150	ILE
1	E	155	LEU
1	E	162	LYS
1	E	167	ILE
1	E	169	THR
1	E	200	ILE
1	E	237	ASP
1	E	284	VAL
1	E	308	HIS
1	E	337	LYS
1	E	346	LYS
1	E	368	ASN
1	E	391	ARG
1	E	400	LYS
1	F	55	SER
1	F	56	LYS
1	F	68	ILE
1	F	85	LEU
1	F	92	GLU
1	F	105	GLU
1	F	115	LYS
1	F	138	GLN
1	F	145	ILE
1	F	146	GLU
1	F	147	LYS
1	F	149	GLU
1	F	155	LEU
1	F	234	ASN
1	F	237	ASP
1	F	284	VAL
1	F	308	HIS
1	F	333	ARG

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Mol	Chain	Res	Type
1	F	337	LYS
1	F	346	LYS
1	F	368	ASN
1	F	391	ARG
1	F	400	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	117	ASN
1	A	154	ASN
1	A	164	ASN
1	A	214	ASN
1	A	230	GLN
1	B	57	ASN
1	B	154	ASN
1	B	164	ASN
1	B	230	GLN
1	C	154	ASN
1	C	214	ASN
1	C	230	GLN
1	D	140	ASN
1	D	154	ASN
1	D	164	ASN
1	D	230	GLN
1	E	57	ASN
1	E	140	ASN
1	E	154	ASN
1	E	164	ASN
1	E	230	GLN
1	F	73	GLN
1	F	154	ASN
1	F	230	GLN

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	UTP	B	501	-	29,30,30	1.64	6 (20%)	43,47,47	1.63	6 (13%)
2	UTP	C	501	-	29,30,30	1.35	5 (17%)	43,47,47	1.53	6 (13%)
2	UTP	D	501	-	29,30,30	1.42	7 (24%)	43,47,47	1.73	6 (13%)
2	UTP	A	501	-	29,30,30	1.76	6 (20%)	43,47,47	1.65	7 (16%)

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	UTP	B	501	-	-	11/22/38/38	0/2/2/2
2	UTP	C	501	-	-	3/22/38/38	0/2/2/2
2	UTP	D	501	-	-	4/22/38/38	0/2/2/2
2	UTP	A	501	-	-	6/22/38/38	0/2/2/2

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	UTP	PA-O3A	4.57	1.64	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	UTP	PB-O3A	4.14	1.64	1.59
2	A	501	UTP	PB-O3B	3.81	1.63	1.59
2	B	501	UTP	PB-O3B	3.81	1.63	1.59
2	B	501	UTP	PB-O3A	3.77	1.63	1.59
2	B	501	UTP	PA-O3A	3.74	1.63	1.59
2	C	501	UTP	PB-O3A	3.16	1.62	1.59
2	A	501	UTP	C4-N3	-2.98	1.33	1.38
2	C	501	UTP	PA-O3A	2.89	1.62	1.59
2	D	501	UTP	PB-O3A	2.87	1.62	1.59
2	D	501	UTP	PB-O3B	2.82	1.62	1.59
2	D	501	UTP	PA-O3A	2.74	1.62	1.59
2	D	501	UTP	C2-N1	2.70	1.42	1.38
2	D	501	UTP	C4-N3	-2.59	1.34	1.38
2	B	501	UTP	C2-N1	2.58	1.42	1.38
2	B	501	UTP	C4-N3	-2.51	1.34	1.38
2	A	501	UTP	C2-N3	-2.45	1.33	1.38
2	C	501	UTP	PB-O3B	2.40	1.62	1.59
2	C	501	UTP	C4-N3	-2.31	1.34	1.38
2	A	501	UTP	C5-C4	-2.17	1.39	1.43
2	B	501	UTP	C5-C4	-2.14	1.39	1.43
2	D	501	UTP	C2-N3	-2.13	1.34	1.38
2	D	501	UTP	C5-C4	-2.10	1.39	1.43
2	C	501	UTP	C6-C5	2.10	1.40	1.35

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	501	UTP	C4-N3-C2	-5.10	120.28	126.61
2	B	501	UTP	C4-N3-C2	-4.98	120.44	126.61
2	C	501	UTP	C4-N3-C2	-4.71	120.77	126.61
2	B	501	UTP	N3-C2-N1	4.49	120.73	114.89
2	D	501	UTP	N3-C2-N1	4.45	120.69	114.89
2	D	501	UTP	C5-C4-N3	4.37	120.92	114.80
2	C	501	UTP	N3-C2-N1	4.33	120.52	114.89
2	A	501	UTP	C4-N3-C2	-4.31	121.26	126.61
2	A	501	UTP	N3-C2-N1	4.24	120.41	114.89
2	A	501	UTP	O4'-C1'-N1	3.93	117.26	108.36
2	B	501	UTP	C5-C4-N3	3.84	120.18	114.80
2	C	501	UTP	C5-C4-N3	3.74	120.04	114.80
2	A	501	UTP	C5-C4-N3	3.71	120.00	114.80
2	D	501	UTP	O4-C4-C5	-3.62	118.93	125.16
2	B	501	UTP	O4-C4-C5	-3.55	119.04	125.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	UTP	O2-C2-N1	-2.93	118.99	122.80
2	D	501	UTP	C3'-C2'-C1'	2.54	106.26	101.46
2	A	501	UTP	O5'-PA-O1A	-2.48	99.11	108.94
2	A	501	UTP	O4-C4-C5	-2.47	120.90	125.16
2	B	501	UTP	O3G-PG-O1G	2.45	116.99	107.80
2	C	501	UTP	O4-C4-C5	-2.39	121.03	125.16
2	D	501	UTP	C1'-N1-C2	2.28	121.69	117.59
2	C	501	UTP	O4'-C1'-N1	2.23	113.42	108.36
2	B	501	UTP	O2-C2-N1	-2.12	120.03	122.80
2	A	501	UTP	O2A-PA-O1A	2.10	122.23	112.44

There are no chirality outliers.

All (24) torsion outliers are listed below:

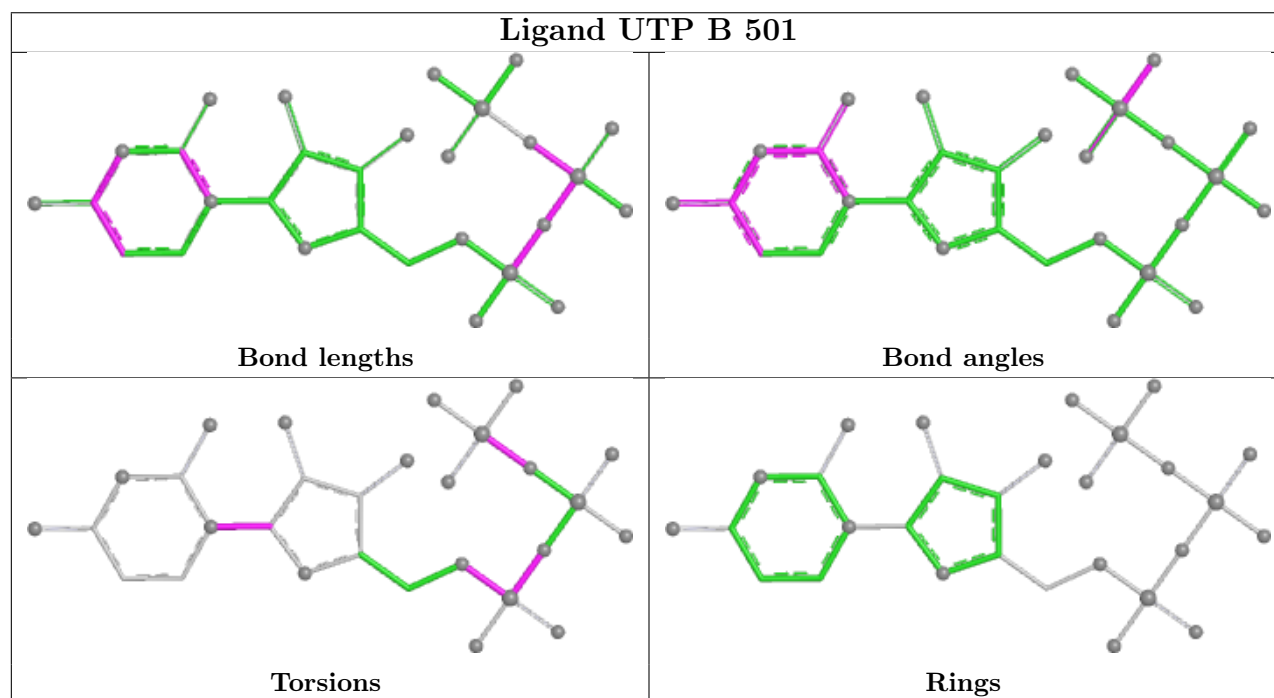
Mol	Chain	Res	Type	Atoms
2	A	501	UTP	C5'-O5'-PA-O3A
2	B	501	UTP	PB-O3A-PA-O5'
2	B	501	UTP	C5'-O5'-PA-O2A
2	B	501	UTP	C5'-O5'-PA-O3A
2	B	501	UTP	PB-O3B-PG-O1G
2	B	501	UTP	PB-O3B-PG-O3G
2	C	501	UTP	C5'-O5'-PA-O3A
2	D	501	UTP	C5'-O5'-PA-O1A
2	D	501	UTP	C5'-O5'-PA-O2A
2	D	501	UTP	C5'-O5'-PA-O3A
2	A	501	UTP	PB-O3B-PG-O2G
2	A	501	UTP	C5'-O5'-PA-O1A
2	C	501	UTP	C5'-O5'-PA-O1A
2	C	501	UTP	C4'-C5'-O5'-PA
2	D	501	UTP	C4'-C5'-O5'-PA
2	B	501	UTP	O4'-C1'-N1-C6
2	A	501	UTP	PG-O3B-PB-O2B
2	B	501	UTP	C2'-C1'-N1-C6
2	A	501	UTP	PB-O3A-PA-O2A
2	A	501	UTP	PG-O3B-PB-O1B
2	B	501	UTP	PB-O3A-PA-O1A
2	B	501	UTP	PB-O3B-PG-O2G
2	B	501	UTP	O4'-C1'-N1-C2
2	B	501	UTP	PB-O3A-PA-O2A

There are no ring outliers.

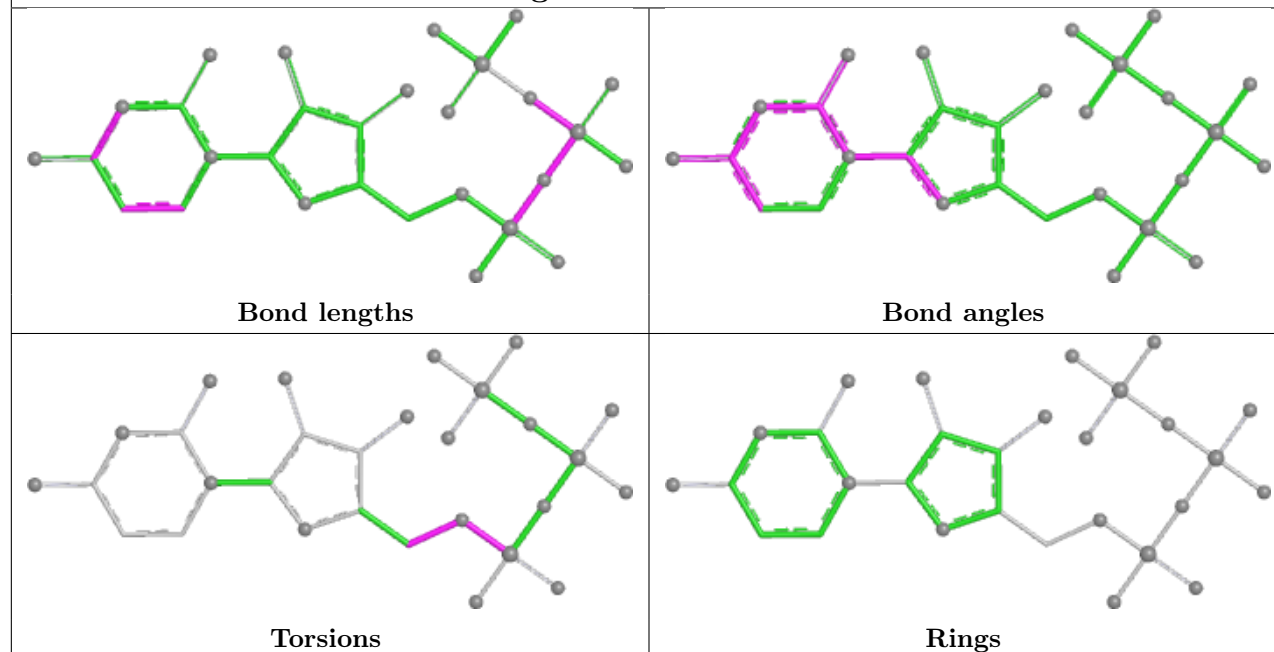
2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	UTP	1	0
2	C	501	UTP	1	0

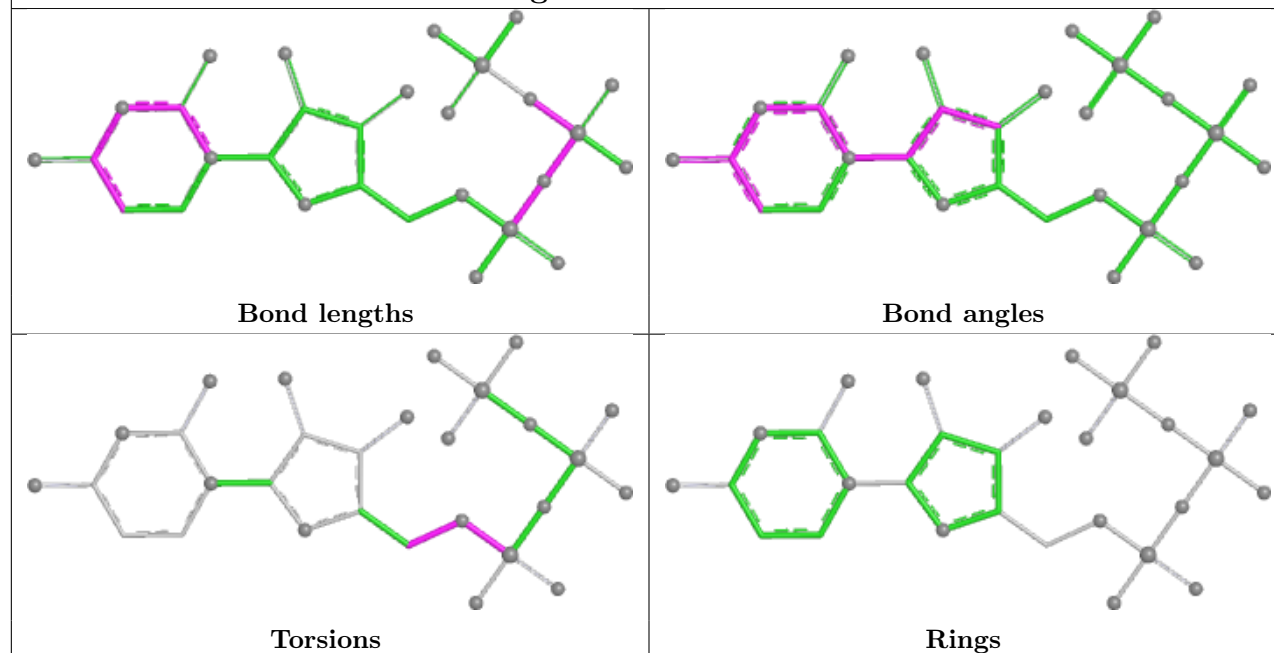
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

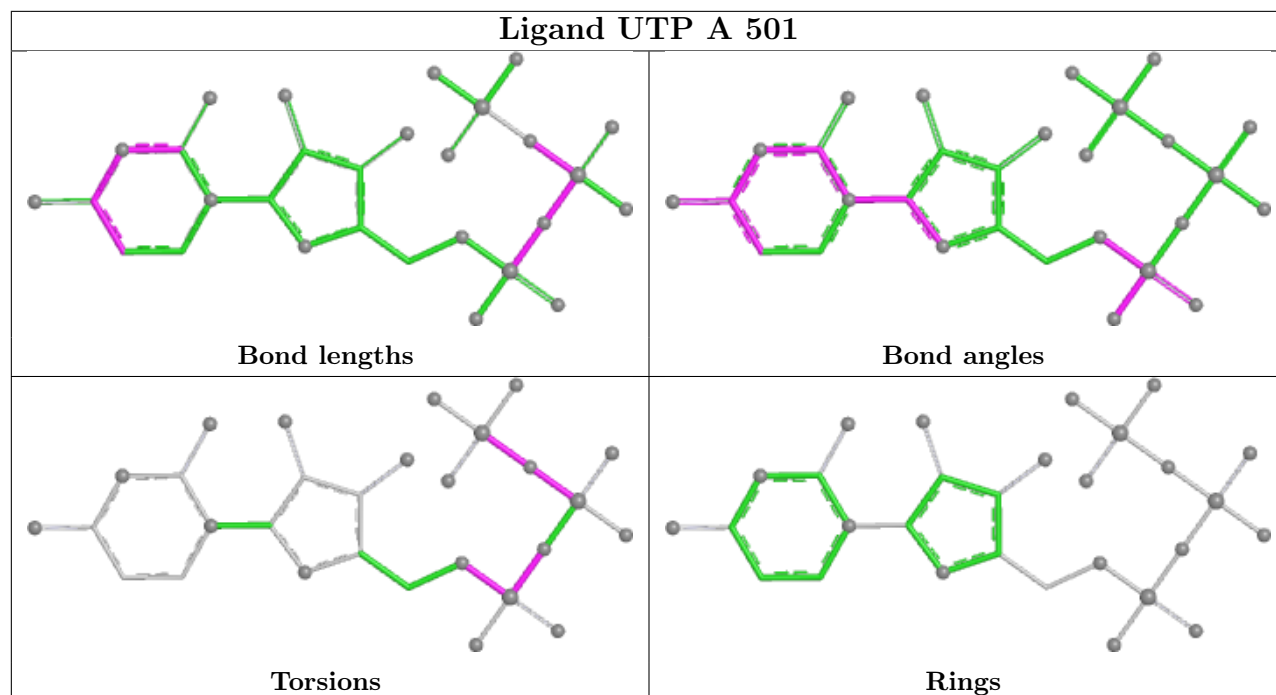


Ligand UTP C 501



Ligand UTP D 501





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	401/409 (98%)	-0.22	6 (1%)	72 64	35, 60, 113, 151	0
1	B	401/409 (98%)	-0.22	2 (0%)	87 83	35, 65, 123, 170	0
1	C	401/409 (98%)	-0.06	6 (1%)	72 64	43, 78, 132, 171	0
1	D	401/409 (98%)	0.03	4 (0%)	79 72	43, 81, 143, 178	0
1	E	369/409 (90%)	0.16	6 (1%)	70 62	49, 86, 138, 171	0
1	F	364/409 (88%)	0.13	7 (1%)	66 58	39, 84, 165, 205	0
All	All	2337/2454 (95%)	-0.04	31 (1%)	75 67	35, 74, 142, 205	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	334	PHE	4.5
1	B	401	VAL	3.9
1	D	401	VAL	3.3
1	C	150	ILE	3.0
1	C	401	VAL	3.0
1	D	96	ILE	2.9
1	A	401	VAL	2.9
1	E	81	GLY	2.8
1	E	209	ILE	2.7
1	A	194	ASP	2.7
1	D	228	PHE	2.6
1	F	345	GLY	2.6
1	E	401	VAL	2.6
1	E	163	LEU	2.5
1	A	150	ILE	2.5
1	A	228	PHE	2.4
1	C	196	ARG	2.4
1	C	87	ALA	2.4
1	E	82	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	F	119	ILE	2.3
1	F	95	ILE	2.3
1	D	107	GLU	2.2
1	F	401	VAL	2.2
1	E	226	LEU	2.2
1	F	78	LYS	2.2
1	B	182	LEU	2.2
1	C	176	ILE	2.1
1	F	83	ALA	2.1
1	F	98	GLY	2.1
1	C	77	ILE	2.1
1	A	174	ILE	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](#)

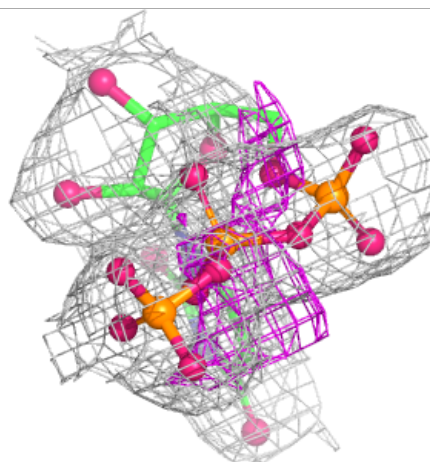
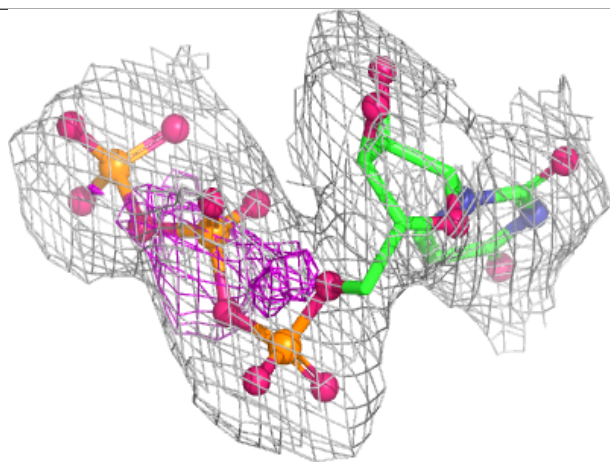
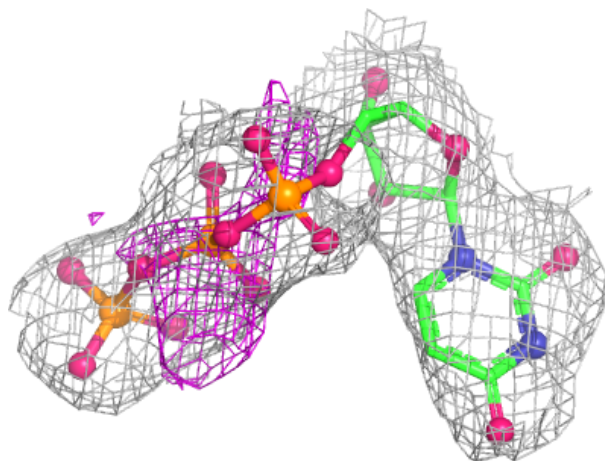
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	UTP	B	501	29/29	0.86	0.09	61,103,144,159	0
2	UTP	C	501	29/29	0.88	0.07	77,98,137,154	0
2	UTP	D	501	29/29	0.92	0.06	65,82,121,128	0
2	UTP	A	501	29/29	0.93	0.07	48,62,79,101	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.

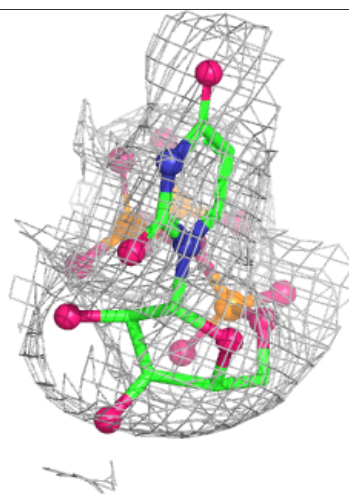
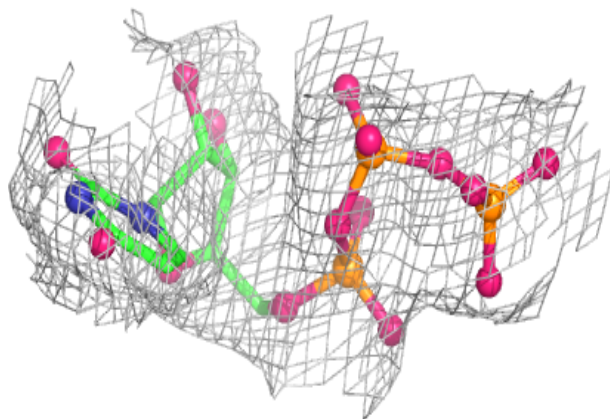
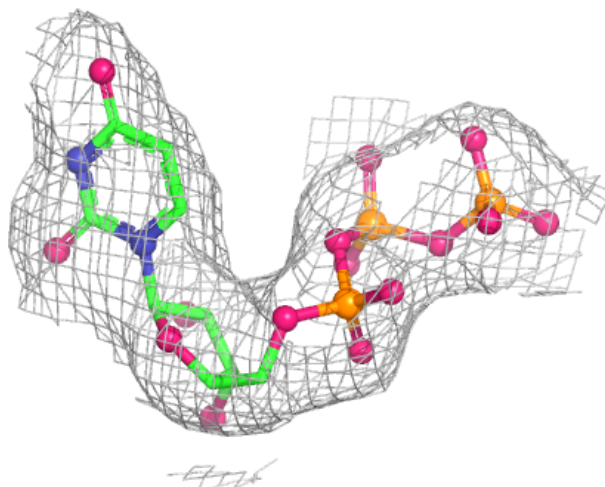
Electron density around UTP B 501:

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



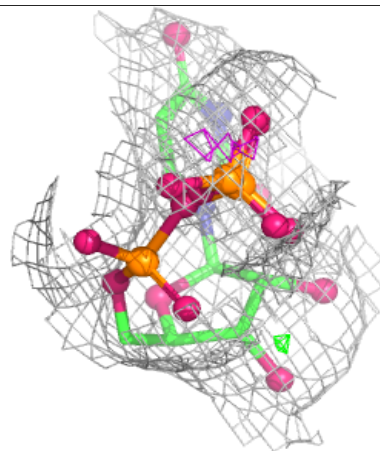
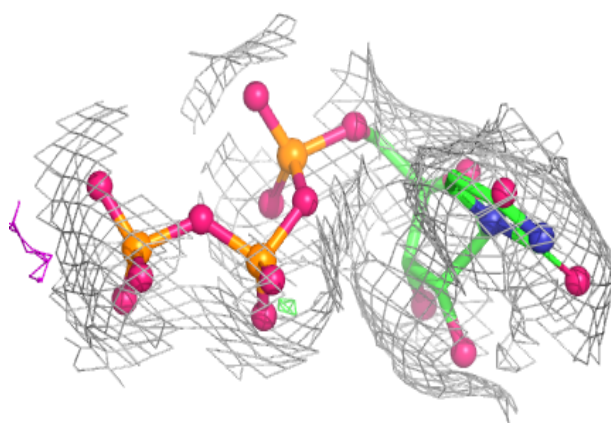
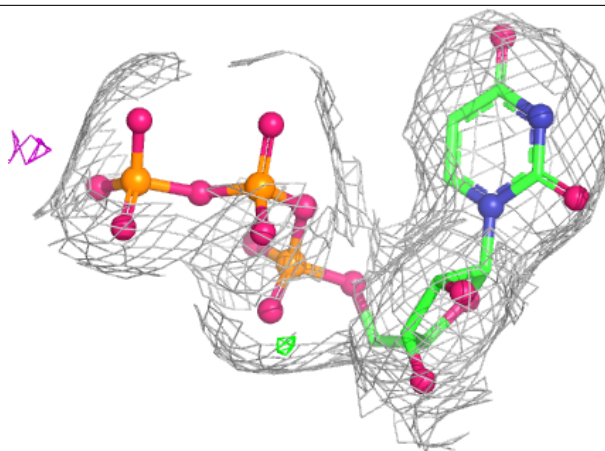
Electron density around UTP C 501:

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



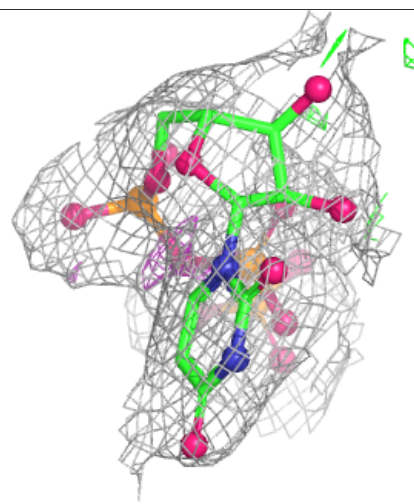
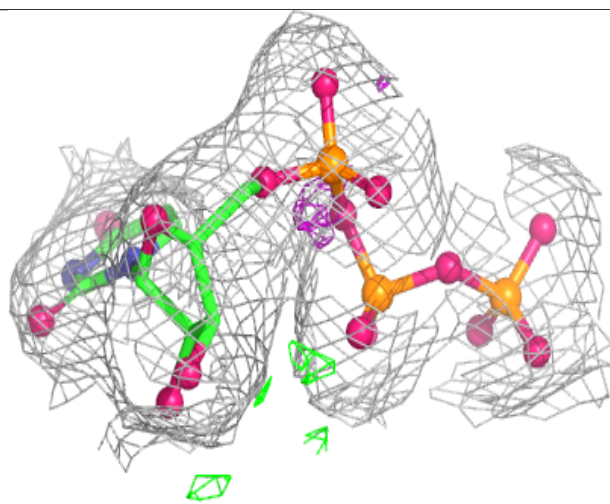
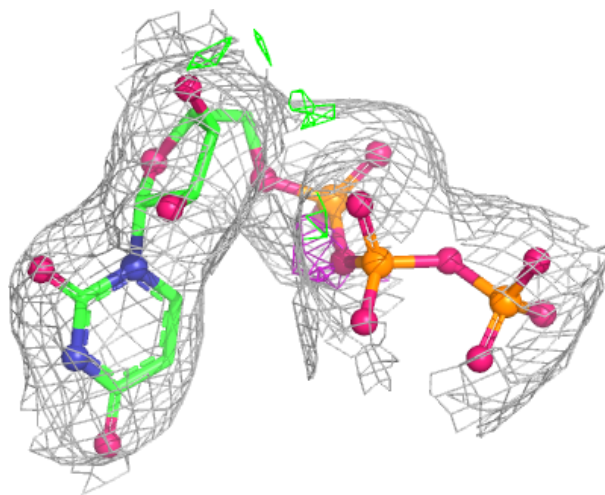
Electron density around UTP D 501:

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



Electron density around UTP A 501:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.