



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 12:06 PM UTC

PDB ID : 8B79 / pdb_00008b79
Title : The crystal structure of M644G variant of DNA Pol Epsilon containing UTP
in the polymerase active site
Authors : Parkash, V.; Johansson, E.
Deposited on : 2022-09-29
Resolution : 2.65 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

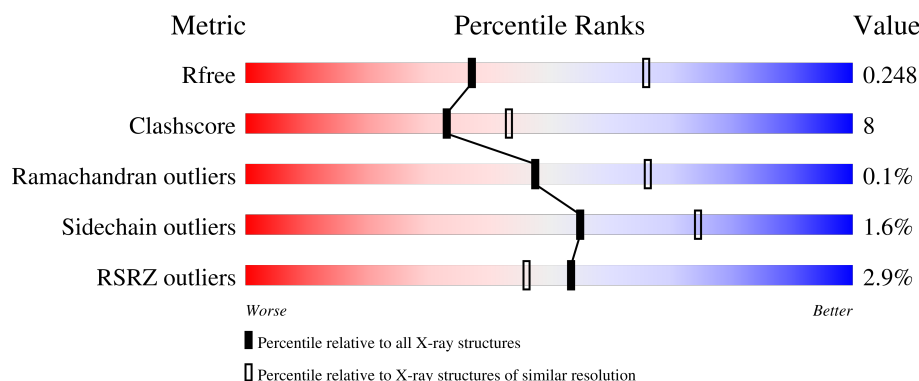
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 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	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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	1191	2% 74% 19% 6%
1	B	1191	3% 76% 17% 6%
2	C	11	73% 27%
2	P	11	9% 91% 9%
3	D	16	69% 25% 6%

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Mol	Chain	Length	Quality of chain
3	T	16	88% 6% 6%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 19091 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 DNA polymerase epsilon catalytic subunit A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1118	Total	C	N	O	S	0	0	0
			9014	5773	1500	1698	43			
1	B	1119	Total	C	N	O	S	0	0	0
			8937	5715	1494	1687	41			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	expression tag	UNP P21951
A	-3	GLY	-	expression tag	UNP P21951
A	-2	ASP	-	expression tag	UNP P21951
A	-1	PRO	-	expression tag	UNP P21951
A	0	HIS	-	expression tag	UNP P21951
A	644	GLY	MET	engineered mutation	UNP P21951
B	-4	GLY	-	expression tag	UNP P21951
B	-3	GLY	-	expression tag	UNP P21951
B	-2	ASP	-	expression tag	UNP P21951
B	-1	PRO	-	expression tag	UNP P21951
B	0	HIS	-	expression tag	UNP P21951
B	644	GLY	MET	engineered mutation	UNP P21951

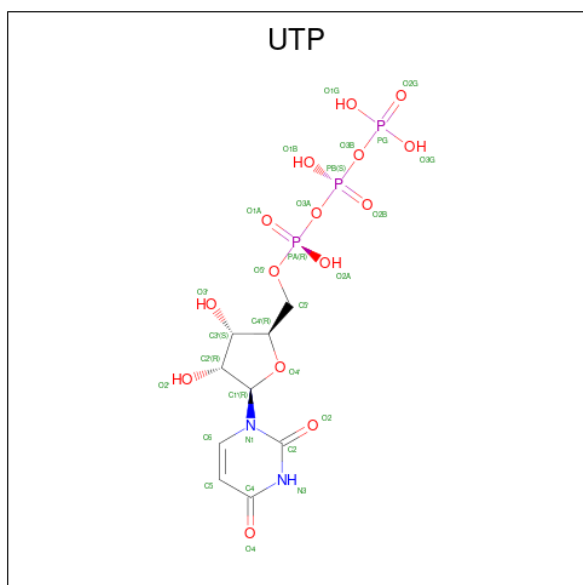
- Molecule 2 is a DNA chain called Primer DNA sequence.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	11	Total	C	N	O	P	0	0	0
			217	106	38	63	10			
2	P	11	Total	C	N	O	P	0	0	0
			221	106	38	66	11			

- Molecule 3 is a DNA chain called Template DNA Sequence.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	15	Total	C	N	O	P	0	0	0
			309	147	57	90	15			
3	T	15	Total	C	N	O	P	0	0	0
			309	147	57	90	15			

- Molecule 4 is URIDINE 5'-TRIPHOSPHATE (CCD ID: UTP) (formula: $C_9H_{15}N_2O_{15}P_3$) (labeled as "Ligand of Interest" by depositor).



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

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	3	Total	Ca	0	0
			3	3		
5	B	3	Total	Ca	0	0
			3	3		

- 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	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		

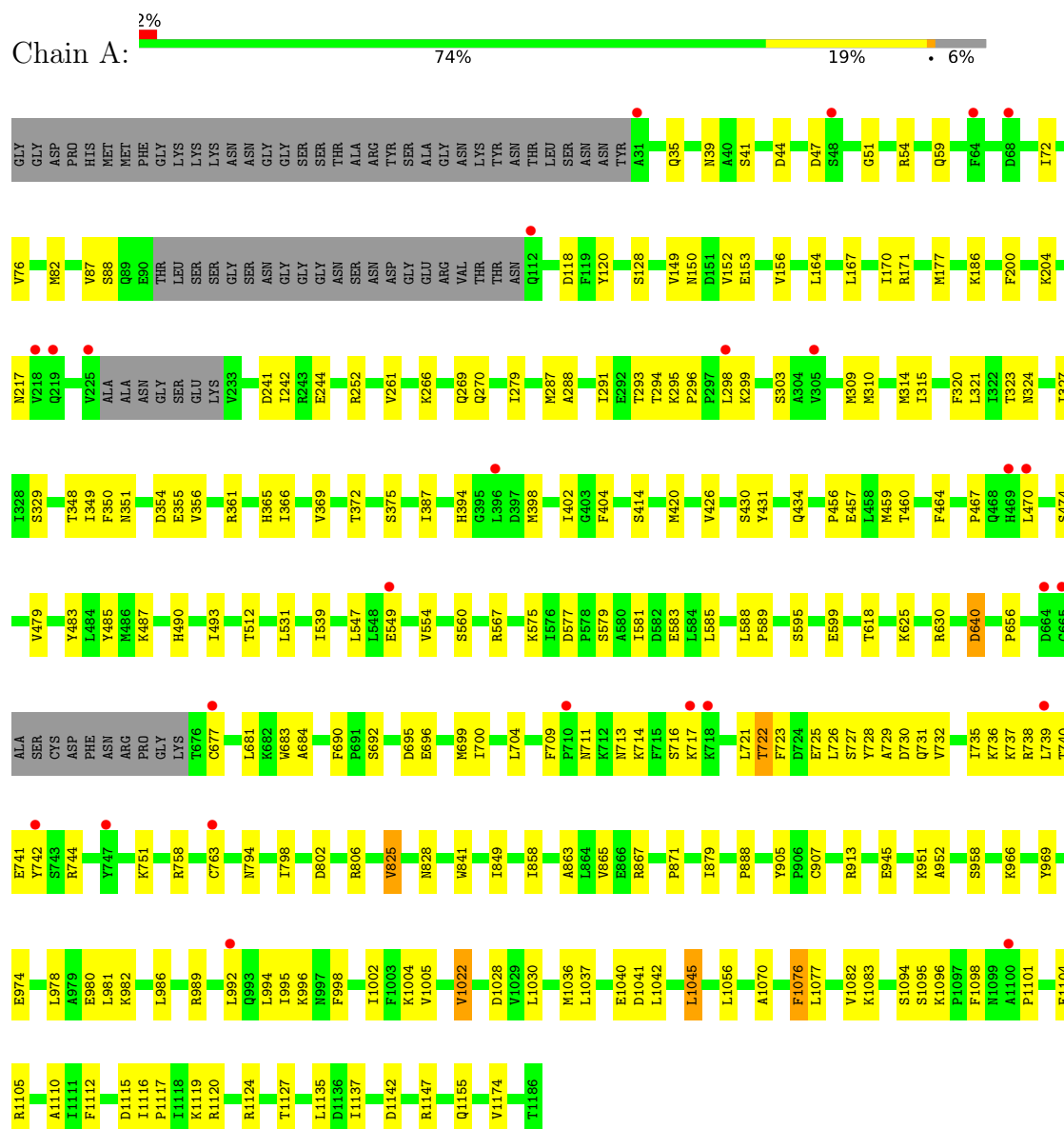
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	2	Total	O	0	0
			2	2		
7	B	6	Total	O	0	0
			6	6		
7	D	2	Total	O	0	0
			2	2		
7	P	1	Total	O	0	0
			1	1		
7	T	1	Total	O	0	0
			1	1		

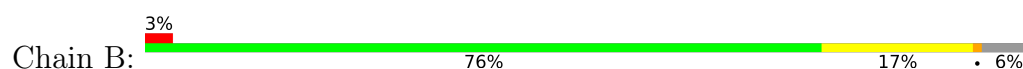
3 Residue-property plots [i](#)

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: DNA polymerase epsilon catalytic subunit A



- Molecule 1: DNA polymerase epsilon catalytic subunit A





- Molecule 3: Template DNA Sequence

Chain T: 88% 6% 6%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	154.40Å 70.70Å 159.40Å 90.00° 112.80° 90.00°	Depositor
Resolution (Å)	43.54 – 2.65 43.54 – 2.65	Depositor EDS
% Data completeness (in resolution range)	99.8 (43.54-2.65) 99.9 (43.54-2.65)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.21 (at 2.65Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.215 , 0.248 0.217 , 0.248	Depositor DCC
R_{free} test set	4636 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	73.4	Xtriage
Anisotropy	0.143	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 42.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	19091	wwPDB-VP
Average B, all atoms (Å ²)	75.0	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 56.67 % 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 2.6480e-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: DOC, CA, ACT, 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	0.27	0/9220	0.47	0/12479
1	B	0.26	0/9140	0.48	1/12386 (0.0%)
2	C	0.27	0/222	0.54	0/341
2	P	0.26	0/226	0.52	0/346
3	D	0.18	0/346	0.37	0/532
3	T	0.18	0/346	0.37	0/532
All	All	0.27	0/19500	0.47	1/26616 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	588	LEU	N-CA-C	-5.31	105.62	113.16

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	9014	0	8824	147	0
1	B	8937	0	8672	139	0
2	C	217	0	123	3	0
2	P	221	0	125	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	309	0	170	3	0
3	T	309	0	170	1	0
4	A	29	0	11	2	0
4	B	29	0	11	1	0
5	A	3	0	0	0	0
5	B	3	0	0	0	0
6	A	4	0	3	0	0
6	B	4	0	3	0	0
7	A	2	0	0	0	0
7	B	6	0	0	0	0
7	D	2	0	0	0	0
7	P	1	0	0	0	0
7	T	1	0	0	0	0
All	All	19091	0	18112	288	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 8.

All (288) 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:639:VAL:HG12	1:B:946:VAL:HG22	1.43	0.97
1:A:700:ILE:HG21	1:A:738:ARG:HB3	1.59	0.82
1:B:160:LEU:HD11	1:B:206:LEU:HD21	1.70	0.73
1:A:986:LEU:HA	1:A:996:LYS:HD3	1.73	0.70
1:A:314:MET:HE1	1:A:479:VAL:HA	1.74	0.69
1:A:547:LEU:HD21	1:A:736:LYS:HE2	1.73	0.69
1:A:1137:ILE:HD12	1:A:1137:ILE:H	1.57	0.69
1:B:792:LYS:HG2	1:B:813:ILE:HD13	1.74	0.68
1:A:1112:PHE:HA	1:A:1119:LYS:HD2	1.76	0.67
1:B:919:THR:HA	1:B:940:ASN:HD22	1.59	0.67
1:B:209:ILE:HG21	1:B:239:ILE:HD12	1.77	0.66
1:B:164:LEU:HD21	1:B:167:LEU:HD23	1.77	0.66
1:A:974:GLU:H	1:A:974:GLU:CD	2.03	0.66
1:B:913:ARG:O	1:B:917:LYS:HG3	1.95	0.66
1:A:736:LYS:O	1:A:740:THR:HG23	1.96	0.66
1:A:298:LEU:HD13	1:A:1110:ALA:HB1	1.77	0.66
1:B:893:PHE:HB2	1:B:901:LEU:HB2	1.77	0.65
1:A:287:MET:HG3	1:A:315:ILE:HG12	1.77	0.65
1:A:583:GLU:OE1	1:A:867:ARG:NH2	2.30	0.64
1:B:329:SER:HB3	1:B:464:PHE:HA	1.78	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:315:ILE:HD13	1:B:369:VAL:HG11	1.79	0.63
1:A:76:VAL:HG22	1:A:266:LYS:HG3	1.80	0.63
1:B:1049:ASN:HD21	1:B:1088:GLN:HB3	1.64	0.63
1:A:315:ILE:HG21	1:A:369:VAL:HG11	1.79	0.63
1:A:690:PHE:HE2	1:A:740:THR:HG22	1.63	0.63
1:A:82:MET:HE3	1:A:261:VAL:HG23	1.80	0.62
1:B:612:GLU:O	1:B:616:GLN:HG3	1.98	0.62
1:B:695:ASP:OD1	1:B:696:GLU:N	2.29	0.62
1:A:398:MET:HG2	1:A:402:ILE:HD11	1.80	0.62
1:A:296:PRO:HB2	1:A:299:LYS:HB2	1.83	0.61
1:A:512:THR:HG21	1:A:825:VAL:HG22	1.81	0.61
1:A:329:SER:HB3	1:A:464:PHE:HA	1.83	0.60
1:B:548:LEU:HD23	1:B:689:PHE:HB3	1.83	0.60
1:A:1056:LEU:HB3	1:A:1082:VAL:HG11	1.82	0.60
1:B:1112:PHE:CG	1:B:1137:ILE:HD11	2.37	0.59
1:B:1092:ILE:HG23	1:B:1109:VAL:HG12	1.84	0.59
1:B:1011:THR:HG23	1:B:1014:GLY:H	1.68	0.59
1:A:287:MET:HE1	1:A:366:ILE:HG13	1.85	0.58
1:B:588:LEU:N	1:B:589:PRO:CD	2.66	0.58
1:A:588:LEU:HB3	1:A:589:PRO:HD3	1.85	0.58
1:A:969:TYR:CZ	1:A:982:LYS:HG3	2.39	0.58
1:B:681:LEU:HB2	1:B:849:ILE:HD13	1.86	0.58
1:A:728:TYR:O	1:A:732:VAL:HG23	2.03	0.58
1:A:310:MET:HE3	1:A:474:SER:HB3	1.85	0.58
1:A:177:MET:HE1	1:A:186:LYS:HB2	1.86	0.57
1:A:350:PHE:CG	1:A:361:ARG:HD3	2.39	0.57
1:B:455:ASP:HB3	1:B:458:LEU:HG	1.85	0.57
1:A:722:THR:HG22	1:A:725:GLU:HG3	1.86	0.57
1:A:989:ARG:HH21	2:P:8:DG:P	2.28	0.57
1:B:664:ASP:O	1:B:665:CYS:HB3	2.04	0.57
1:A:35:GLN:HE21	1:A:39:ASN:HD21	1.51	0.56
1:B:342:GLU:CD	1:B:342:GLU:H	2.13	0.56
1:B:291:ILE:HG13	1:B:383:ASP:OD2	2.05	0.56
1:B:1124:ARG:NH1	1:B:1132:LEU:O	2.38	0.56
1:A:951:LYS:HB2	1:A:974:GLU:HA	1.87	0.56
1:B:924:GLN:NE2	1:B:936:THR:OG1	2.35	0.56
1:B:1112:PHE:CE1	1:B:1123:LEU:HD21	2.39	0.56
1:A:577:ASP:OD1	1:A:579:SER:OG	2.22	0.56
1:A:1037:LEU:HD12	1:A:1042:LEU:HD13	1.86	0.55
1:A:711:ASN:ND2	1:A:716:SER:HB3	2.22	0.55
1:B:1079:GLU:O	1:B:1079:GLU:OE1	2.25	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:700:ILE:CG2	1:A:738:ARG:HB3	2.34	0.54
1:A:1116:ILE:N	1:A:1117:PRO:HD2	2.22	0.54
1:A:863:ALA:O	1:A:867:ARG:HG3	2.07	0.54
1:B:613:ILE:HD12	1:B:891:TYR:HB3	1.90	0.54
1:B:621:LEU:HD21	1:B:888:PRO:HG3	1.90	0.54
1:A:1120:ARG:HB2	1:A:1135:LEU:HD11	1.89	0.54
1:B:146:GLU:HG2	1:B:187:THR:HB	1.90	0.54
1:B:910:LEU:O	1:B:914:VAL:HG23	2.07	0.54
1:A:539:ILE:HD12	1:A:728:TYR:CD2	2.43	0.54
1:B:728:TYR:O	1:B:732:VAL:HG23	2.08	0.54
1:B:387:ILE:O	1:B:391:SER:OG	2.21	0.53
1:B:321:LEU:HB3	1:B:349:ILE:HD12	1.91	0.53
1:B:609:ASN:HD21	1:B:893:PHE:HA	1.72	0.53
1:B:693:LYS:CD	1:B:694:MET:H	2.20	0.53
1:A:865:VAL:HG12	1:A:871:PRO:HD3	1.90	0.53
1:A:244:GLU:OE1	1:A:252:ARG:NH2	2.41	0.53
1:A:1056:LEU:HD13	1:A:1070:ALA:HB3	1.90	0.53
1:A:980:GLU:HB3	1:A:982:LYS:HE2	1.90	0.53
1:B:924:GLN:HE21	1:B:936:THR:HG1	1.56	0.53
1:A:87:VAL:HG22	1:A:88:SER:O	2.09	0.52
1:A:581:ILE:HG21	1:A:625:LYS:HB2	1.90	0.52
1:A:1096:LYS:HB3	1:A:1127:THR:OG1	2.09	0.52
1:A:828:ASN:HD22	4:A:1201:UTP:H2'	1.75	0.52
1:A:695:ASP:OD1	1:A:696:GLU:N	2.41	0.52
1:A:585:LEU:HD21	1:A:618:THR:HG23	1.90	0.52
1:A:713:ASN:O	1:A:714:LYS:C	2.53	0.52
1:B:685:TRP:HB3	1:B:756:VAL:HG22	1.92	0.52
1:B:1024:ASN:OD1	1:B:1173:ARG:HD3	2.09	0.52
1:B:552:THR:OG1	1:B:553:TYR:N	2.43	0.51
1:B:340:LYS:HB3	1:B:342:GLU:OE1	2.09	0.51
1:A:1120:ARG:HD3	1:A:1124:ARG:CZ	2.41	0.51
1:B:861:ALA:O	1:B:865:VAL:HG23	2.09	0.51
1:B:338:THR:HG23	1:B:344:PRO:HA	1.92	0.51
1:A:695:ASP:O	1:A:699:MET:HG3	2.11	0.51
1:A:727:SER:O	1:A:731:GLN:HG3	2.10	0.51
1:B:571:LYS:HG2	1:B:633:LEU:HD23	1.93	0.51
1:A:992:LEU:HD23	1:A:995:ILE:HG12	1.93	0.51
1:B:354:ASP:OD1	1:B:354:ASP:N	2.42	0.51
1:A:270:GLN:HA	1:A:270:GLN:OE1	2.10	0.50
1:B:907:CYS:SG	1:B:946:VAL:HG23	2.51	0.50
1:A:321:LEU:HB3	1:A:349:ILE:HD13	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:320:PHE:HA	1:A:348:THR:O	2.12	0.50
1:A:1096:LYS:HG3	1:A:1098:PHE:CE1	2.46	0.50
1:B:588:LEU:N	1:B:589:PRO:HD2	2.27	0.49
1:B:244:GLU:OE1	1:B:252:ARG:NH2	2.45	0.49
2:C:10:DT:H2''	2:C:11:DOC:H5'	1.95	0.49
1:B:693:LYS:HD3	1:B:694:MET:H	1.78	0.49
1:A:150:ASN:ND2	1:A:153:GLU:OE2	2.46	0.49
1:A:656:PRO:HG2	1:A:841:TRP:HB3	1.93	0.49
1:B:355:GLU:HB3	1:B:394:HIS:HE2	1.78	0.49
1:B:568:SER:O	1:B:633:LEU:HD13	2.13	0.49
1:A:794:ASN:O	1:A:798:ILE:HG13	2.12	0.49
1:B:588:LEU:HB3	1:B:589:PRO:HD3	1.95	0.49
1:B:746:VAL:HG13	1:B:747:TYR:CD2	2.47	0.49
1:B:789:LYS:HD3	1:B:789:LYS:C	2.38	0.49
1:B:32:LEU:O	1:B:36:GLN:HG3	2.13	0.48
1:B:207:ARG:O	1:B:211:GLN:HG3	2.14	0.48
1:B:621:LEU:O	1:B:624:LEU:HB2	2.13	0.48
1:B:696:GLU:O	1:B:700:ILE:HG12	2.13	0.48
1:B:512:THR:HG21	1:B:825:VAL:HG12	1.95	0.48
1:B:68:ASP:O	1:B:72:ILE:HG13	2.14	0.48
1:B:645:TYR:O	1:B:649:MET:HG3	2.14	0.48
1:B:956:PRO:HG2	1:B:965:ILE:HG21	1.96	0.48
1:B:928:ASP:HB3	1:B:933:ILE:HB	1.96	0.47
1:A:456:PRO:HA	1:A:459:MET:HG2	1.96	0.47
1:B:393:ILE:HG22	1:B:394:HIS:ND1	2.28	0.47
1:B:591:ALA:HA	1:B:912:TYR:CD1	2.49	0.47
1:B:383:ASP:O	1:B:387:ILE:HG13	2.14	0.47
1:A:47:ASP:OD2	1:A:120:TYR:OH	2.30	0.47
1:A:802:ASP:O	1:A:806:ARG:HG3	2.15	0.47
1:B:462:TYR:CE2	1:B:466:LYS:HD2	2.49	0.47
1:A:170:ILE:HD12	1:A:171:ARG:H	1.79	0.47
1:A:365:HIS:O	1:A:369:VAL:HG23	2.14	0.47
1:A:737:LYS:O	1:A:741:GLU:HG3	2.14	0.47
1:A:1041:ASP:O	1:A:1045:LEU:HB2	2.14	0.47
1:B:61:ASN:OD1	1:B:61:ASN:N	2.43	0.47
1:B:845:GLU:O	1:B:849:ILE:HG13	2.15	0.47
1:A:729:ALA:O	1:A:730:ASP:C	2.58	0.47
1:A:329:SER:OG	1:A:467:PRO:HG3	2.15	0.47
1:A:709:PHE:CD2	1:A:726:LEU:HD11	2.50	0.47
1:B:722:THR:HG22	1:B:725:GLU:OE2	2.15	0.47
1:B:434:GLN:HB2	2:C:9:DT:OP2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:314:MET:CE	1:A:479:VAL:HA	2.43	0.46
1:B:1025:ARG:HG3	1:B:1026:TRP:N	2.30	0.46
1:B:562:GLU:OE1	1:B:870:ARG:NH2	2.48	0.46
1:B:778:ARG:NH1	1:B:779:ASP:OD1	2.48	0.46
1:B:1180:LEU:O	1:B:1184:ILE:HG13	2.14	0.46
1:B:309:MET:HE3	1:B:470:LEU:HD21	1.97	0.46
1:A:858:ILE:HD11	1:A:879:ILE:HG13	1.97	0.46
1:A:727:SER:C	1:A:729:ALA:N	2.71	0.46
1:A:732:VAL:O	1:A:736:LYS:HG2	2.15	0.46
1:A:200:PHE:CZ	1:A:204:LYS:HE3	2.51	0.46
1:A:549:GLU:OE1	1:A:751:LYS:HE2	2.15	0.46
1:B:235:ALA:HA	1:B:238:LEU:HD23	1.98	0.46
1:B:458:LEU:C	1:B:461:PRO:HD2	2.40	0.46
1:B:1112:PHE:CD1	1:B:1137:ILE:HD11	2.51	0.46
1:A:828:ASN:ND2	4:A:1201:UTP:H2'	2.31	0.45
1:A:279:ILE:HG22	1:A:279:ILE:O	2.15	0.45
1:B:397:ASP:O	1:B:401:GLU:HG2	2.16	0.45
1:A:59:GLN:HA	1:A:269:GLN:OE1	2.17	0.45
1:B:736:LYS:O	1:B:740:THR:HG23	2.16	0.45
1:B:1022:VAL:O	1:B:1026:TRP:HD1	1.99	0.45
1:A:241:ASP:OD1	1:A:242:ILE:N	2.48	0.45
1:A:420:MET:HE1	1:A:493:ILE:HG21	1.98	0.45
1:A:681:LEU:HB2	1:A:849:ILE:HD13	1.98	0.45
1:B:57:PRO:HG3	1:B:267:VAL:HG12	1.99	0.45
1:B:310:MET:HE2	1:B:310:MET:HB3	1.78	0.45
1:A:295:LYS:NZ	1:A:457:GLU:OE2	2.46	0.45
1:A:1028:ASP:HB3	1:A:1036:MET:HE1	1.99	0.45
1:A:981:LEU:HD21	1:A:986:LEU:HD23	1.99	0.44
1:A:696:GLU:HG2	1:A:742:TYR:OH	2.17	0.44
1:A:1115:ASP:O	1:A:1116:ILE:C	2.60	0.44
1:B:485:TYR:CE2	1:B:490:HIS:HB2	2.53	0.44
1:B:640:ASP:O	1:B:945:GLU:HG2	2.17	0.44
1:B:806:ARG:O	1:B:810:LYS:HG3	2.17	0.44
1:B:906:PRO:HA	1:B:909:MET:SD	2.57	0.44
1:A:244:GLU:HG2	1:A:531:LEU:HB2	1.98	0.44
1:A:858:ILE:HD12	1:A:858:ILE:HA	1.87	0.44
1:B:356:VAL:O	1:B:360:GLN:HG3	2.17	0.44
1:A:716:SER:O	1:A:717:LYS:C	2.61	0.44
1:A:1056:LEU:CB	1:A:1082:VAL:HG11	2.47	0.44
1:A:324:ASN:OD1	1:A:355:GLU:N	2.50	0.44
1:A:723:PHE:CZ	1:A:731:GLN:HB3	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:260:ARG:HB2	1:B:263:LYS:HD2	2.00	0.44
1:A:690:PHE:CE2	1:A:740:THR:HG22	2.50	0.44
1:B:433:PRO:HG2	1:B:436:SER:HB2	2.00	0.44
1:A:164:LEU:HD21	1:A:167:LEU:HD23	1.99	0.44
1:A:560:SER:HB2	1:A:966:LYS:HG3	1.99	0.44
1:A:1094:SER:O	1:A:1105:ARG:HD2	2.18	0.43
1:B:627:ASN:HB2	1:B:630:ARG:NH1	2.33	0.43
1:A:704:LEU:HD12	1:A:709:PHE:HZ	1.83	0.43
1:A:44:ASP:OD1	1:A:54:ARG:NH2	2.50	0.43
1:A:723:PHE:O	1:A:726:LEU:HB2	2.18	0.43
1:A:323:THR:OG1	1:A:351:ASN:HA	2.18	0.43
1:B:420:MET:HA	1:B:420:MET:HE2	2.00	0.43
1:B:590:GLU:HA	1:B:593:LYS:HD2	2.00	0.43
1:A:426:VAL:HA	1:A:430:SER:HB3	1.99	0.43
1:A:677:CYS:O	1:A:763:CYS:HA	2.17	0.43
1:B:664:ASP:O	1:B:665:CYS:CB	2.66	0.43
1:B:1094:SER:O	1:B:1105:ARG:HD2	2.18	0.43
1:B:1124:ARG:HG2	1:B:1124:ARG:HH11	1.84	0.43
1:A:640:ASP:O	1:A:945:GLU:HG2	2.18	0.43
1:A:731:GLN:O	1:A:735:ILE:HG13	2.19	0.43
1:A:1076:PHE:CE2	1:A:1077:LEU:HG	2.54	0.43
1:B:194:VAL:HG12	1:B:198:GLN:HG3	2.01	0.43
1:B:1112:PHE:HE1	1:B:1123:LEU:HD21	1.84	0.43
1:A:167:LEU:HD12	1:A:167:LEU:C	2.44	0.43
1:B:722:THR:OG1	1:B:723:PHE:N	2.51	0.43
1:B:722:THR:HG23	1:B:725:GLU:H	1.84	0.43
1:A:404:PHE:CZ	1:A:414:SER:HB3	2.53	0.43
1:B:289:PHE:CG	1:B:362:PHE:HZ	2.37	0.43
1:B:548:LEU:HD23	1:B:548:LEU:HA	1.89	0.43
1:A:888:PRO:HB2	1:A:905:TYR:CD2	2.53	0.43
1:B:361:ARG:HE	1:B:361:ARG:HB2	1.57	0.42
1:B:36:GLN:OE1	1:B:86:LEU:HD22	2.19	0.42
1:B:446:LYS:O	1:B:488:TYR:OH	2.32	0.42
1:B:521:MET:HG2	1:B:531:LEU:HD11	2.01	0.42
1:B:951:LYS:HB2	1:B:974:GLU:HA	2.01	0.42
1:A:291:ILE:HG21	1:A:387:ILE:HD11	2.00	0.42
1:B:731:GLN:O	1:B:735:ILE:HD12	2.19	0.42
1:A:692:SER:OG	1:A:739:LEU:HD11	2.19	0.42
1:A:978:LEU:HD21	1:A:981:LEU:HB2	2.02	0.42
1:B:966:LYS:HD3	3:D:9:DC:H5"	2.01	0.42
1:B:1013:GLU:OE1	1:B:1013:GLU:N	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:11:DC:H2"	3:D:12:DG:C8	2.54	0.42
1:A:294:THR:HG22	1:A:460:THR:HG23	2.02	0.42
1:A:751:LYS:HB2	1:A:751:LYS:HE3	1.82	0.42
1:A:998:PHE:CZ	1:A:1002:ILE:HG13	2.55	0.42
1:A:483:TYR:O	1:A:487:LYS:HB2	2.20	0.42
1:A:288:ALA:HA	1:A:375:SER:O	2.20	0.42
1:A:711:ASN:HD21	1:A:716:SER:HB3	1.84	0.42
1:A:1005:VAL:HG21	1:A:1022:VAL:HG21	2.00	0.42
1:B:583:GLU:OE1	1:B:867:ARG:NH2	2.53	0.42
1:B:1183:LYS:HA	1:B:1183:LYS:HD2	1.78	0.42
1:A:293:THR:C	1:A:309:MET:HE2	2.45	0.42
1:A:310:MET:HE1	1:A:470:LEU:C	2.45	0.42
1:A:683:TRP:CZ2	1:A:758:ARG:HD2	2.55	0.42
1:B:287:MET:O	1:B:374:ILE:HA	2.19	0.42
1:B:517:GLU:OE1	1:B:833:TYR:OH	2.28	0.42
1:B:982:LYS:HB3	2:C:10:DT:H5"	2.02	0.42
1:A:356:VAL:HG22	1:A:394:HIS:HB3	2.02	0.41
1:B:592:LEU:HD23	1:B:592:LEU:HA	1.77	0.41
1:A:171:ARG:HA	1:A:186:LYS:O	2.20	0.41
1:B:322:ILE:HD13	1:B:361:ARG:HG2	2.02	0.41
1:B:656:PRO:HG2	1:B:841:TRP:HB3	2.03	0.41
1:B:728:TYR:O	1:B:731:GLN:HB2	2.19	0.41
1:B:981:LEU:HD21	1:B:986:LEU:HD23	2.01	0.41
1:B:1103:THR:HG23	3:D:14:DT:OP1	2.20	0.41
1:A:354:ASP:OD1	1:A:354:ASP:N	2.53	0.41
1:B:179:ASN:HB3	1:B:182:LEU:HD12	2.01	0.41
1:B:462:TYR:CZ	1:B:466:LYS:HD2	2.55	0.41
1:A:51:GLY:O	1:A:128:SER:OG	2.38	0.41
1:A:994:LEU:HD21	1:A:1030:LEU:HD11	2.02	0.41
1:A:204:LYS:HA	1:A:204:LYS:HD3	1.74	0.41
1:A:287:MET:HE1	1:A:366:ILE:CG1	2.51	0.41
1:B:324:ASN:ND2	1:B:327:ILE:HG13	2.35	0.41
1:B:625:LYS:O	1:B:625:LYS:HG3	2.19	0.41
1:B:787:LEU:HD23	1:B:787:LEU:HA	1.92	0.41
1:A:554:VAL:HB	1:A:684:ALA:HB3	2.03	0.41
1:B:222:ILE:H	1:B:222:ILE:HG13	1.66	0.41
1:B:651:THR:HG23	1:B:940:ASN:HA	2.02	0.41
1:A:149:VAL:O	1:A:153:GLU:HB2	2.20	0.41
1:B:562:GLU:CD	1:B:870:ARG:HH21	2.27	0.41
1:A:252:ARG:HA	1:A:252:ARG:HD3	1.80	0.41
1:A:1095:SER:HA	1:A:1142:ASP:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:217:ASN:HD22	1:A:217:ASN:HA	1.71	0.41
1:A:567:ARG:HG2	1:A:952:ALA:HB2	2.01	0.41
1:A:744:ARG:NH1	3:T:16:DA:H4'	2.36	0.41
1:A:1155:GLN:HG3	1:A:1174:VAL:CG1	2.51	0.41
1:B:559:GLU:HG3	1:B:874:LEU:HD22	2.02	0.41
1:B:1049:ASN:HD22	1:B:1090:LYS:HG2	1.86	0.41
1:B:1182:ARG:HG3	1:B:1182:ARG:HH11	1.86	0.41
1:B:496:LEU:HG	1:B:500:ILE:HD12	2.02	0.41
1:A:152:VAL:O	1:A:156:VAL:HG22	2.21	0.40
1:A:1030:LEU:HD13	1:A:1147:ARG:HA	2.03	0.40
1:A:1101:PRO:HD2	1:A:1104:GLU:OE2	2.20	0.40
1:B:810:LYS:O	1:B:814:VAL:HG23	2.21	0.40
1:A:485:TYR:CE2	1:A:490:HIS:HB2	2.56	0.40
1:A:1112:PHE:CD1	1:A:1137:ILE:HG13	2.55	0.40
1:B:539:ILE:H	1:B:539:ILE:HG13	1.61	0.40
1:B:595:SER:OG	1:B:903:LEU:HD23	2.22	0.40
1:A:72:ILE:O	1:A:266:LYS:HE2	2.21	0.40
1:A:595:SER:HA	1:A:599:GLU:HB2	2.03	0.40
1:A:727:SER:C	1:A:729:ALA:H	2.29	0.40
1:B:324:ASN:O	1:B:328:ILE:HG12	2.22	0.40
1:B:645:TYR:HB2	4:B:1201:UTP:O2'	2.22	0.40
1:A:82:MET:HA	1:A:118:ASP:O	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	1110/1191 (93%)	1067 (96%)	43 (4%)	0	100	100
1	B	1111/1191 (93%)	1064 (96%)	45 (4%)	2 (0%)	43	60
All	All	2221/2382 (93%)	2131 (96%)	88 (4%)	2 (0%)	48	66

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	589	PRO
1	B	555	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	985/1066 (92%)	964 (98%)	21 (2%)	47	69
1	B	966/1066 (91%)	956 (99%)	10 (1%)	68	81
All	All	1951/2132 (92%)	1920 (98%)	31 (2%)	55	74

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	SER
1	A	303	SER
1	A	327	ILE
1	A	372	THR
1	A	431	TYR
1	A	434	GLN
1	A	575	LYS
1	A	630	ARG
1	A	640	ASP
1	A	721	LEU
1	A	722	THR
1	A	825	VAL
1	A	907	CYS
1	A	913	ARG
1	A	958	SER
1	A	1004	LYS
1	A	1022	VAL
1	A	1040	GLU
1	A	1045	LEU
1	A	1076	PHE
1	A	1083	LYS

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Mol	Chain	Res	Type
1	B	310	MET
1	B	327	ILE
1	B	358	LEU
1	B	361	ARG
1	B	431	TYR
1	B	640	ASP
1	B	756	VAL
1	B	1025	ARG
1	B	1076	PHE
1	B	1109	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	GLN
1	A	150	ASN
1	A	168	GLN
1	A	185	GLN
1	A	197	ASN
1	A	213	ASN
1	A	217	ASN
1	A	324	ASN
1	A	388	HIS
1	A	572	ASN
1	A	619	GLN
1	A	638	HIS
1	A	706	ASN
1	A	767	ASN
1	A	828	ASN
1	A	973	ASN
1	A	997	ASN
1	A	1062	GLN
1	B	217	ASN
1	B	237	HIS
1	B	434	GLN
1	B	523	GLN
1	B	616	GLN
1	B	628	ASN
1	B	915	HIS
1	B	924	GLN
1	B	940	ASN
1	B	973	ASN

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Mol	Chain	Res	Type
1	B	1049	ASN
1	B	1088	GLN
1	B	1176	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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	DOC	C	11	3,2	16,19,20	0.40	0	20,26,29	0.29	0
2	DOC	P	11	3,2	16,19,20	0.40	0	20,26,29	0.34	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	DOC	C	11	3,2	-	2/7/18/19	0/2/2/2
2	DOC	P	11	3,2	-	2/7/18/19	0/2/2/2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	11	DOC	C3'-C4'-C5'-O5'
2	C	11	DOC	O4'-C4'-C5'-O5'
2	P	11	DOC	O4'-C4'-C5'-O5'
2	P	11	DOC	C3'-C4'-C5'-O5'

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	11	DOC	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 6 are monoatomic - leaving 4 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
6	ACT	A	1205	5	3,3,3	1.39	1 (33%)	3,3,3	1.36	0
4	UTP	A	1201	5	29,30,30	0.55	0	43,47,47	0.57	0
4	UTP	B	1201	5	29,30,30	0.55	0	43,47,47	0.57	0
6	ACT	B	1205	5	3,3,3	1.38	0	3,3,3	1.36	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	UTP	A	1201	5	-	4/22/38/38	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	UTP	B	1201	5	-	3/22/38/38	0/2/2/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	1205	ACT	CH3-C	2.02	1.57	1.49

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

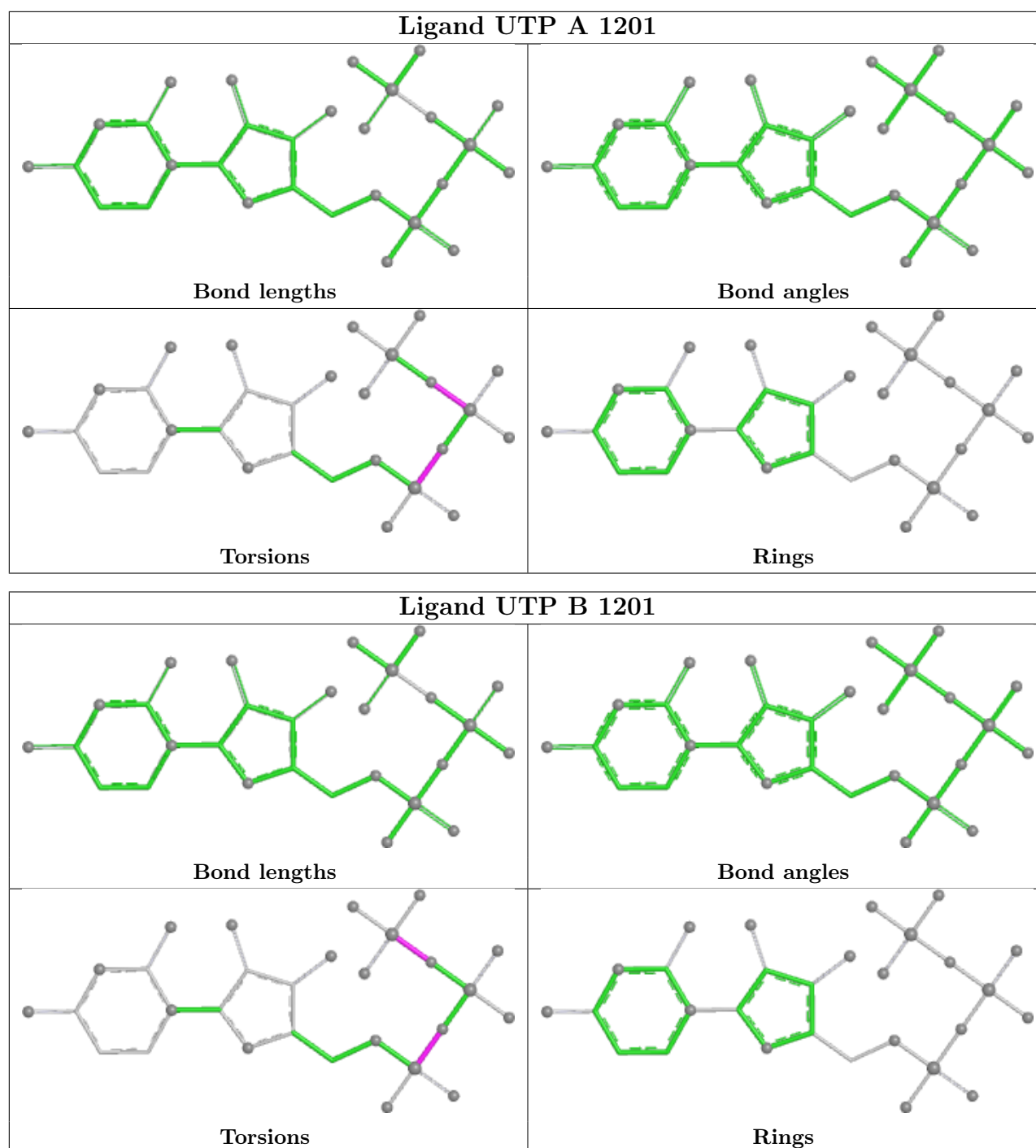
Mol	Chain	Res	Type	Atoms
4	B	1201	UTP	PB-O3B-PG-O3G
4	A	1201	UTP	PG-O3B-PB-O1B
4	B	1201	UTP	PB-O3A-PA-O1A
4	A	1201	UTP	PB-O3A-PA-O2A
4	A	1201	UTP	PB-O3A-PA-O1A
4	B	1201	UTP	PB-O3A-PA-O2A
4	A	1201	UTP	PG-O3B-PB-O2B

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1201	UTP	2	0
4	B	1201	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	1118/1191 (93%)	0.30	26 (2%) 61 54	43, 75, 103, 141	0
1	B	1119/1191 (93%)	0.35	40 (3%) 46 38	45, 76, 110, 139	0
2	C	10/11 (90%)	-0.31	0 100 100	52, 61, 81, 81	0
2	P	10/11 (90%)	-0.02	1 (10%) 12 10	51, 60, 73, 80	1 (10%)
3	D	15/16 (93%)	-0.19	0 100 100	48, 63, 90, 109	0
3	T	15/16 (93%)	-0.19	0 100 100	48, 60, 88, 113	0
All	All	2287/2436 (93%)	0.31	67 (2%) 53 46	43, 75, 107, 141	1 (0%)

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	665	CYS	6.2
1	A	665	CYS	5.3
1	B	677	CYS	4.2
1	B	29	ASN	4.0
1	B	225	VAL	3.7
1	A	677	CYS	3.6
1	B	664	ASP	3.4
1	A	31	ALA	3.3
1	B	220	ARG	3.2
1	B	596	VAL	3.2
1	A	549	GLU	3.2
1	B	742	TYR	3.2
1	A	225	VAL	3.1
1	B	710	PRO	3.1
1	B	690	PHE	3.1
1	B	602	SER	3.0
1	B	595	SER	3.0
1	B	362	PHE	3.0
1	B	747	TYR	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	219	GLN	2.9
1	A	710	PRO	2.9
1	A	992	LEU	2.7
1	B	548	LEU	2.7
1	B	218	VAL	2.7
1	B	1132	LEU	2.7
1	B	663	ARG	2.7
1	A	742	TYR	2.6
1	A	739	LEU	2.6
1	A	763	CYS	2.6
2	P	1	DT	2.6
1	A	64	PHE	2.6
1	A	469	HIS	2.5
1	B	64	PHE	2.5
1	A	664	ASP	2.5
1	B	782	TYR	2.5
1	A	218	VAL	2.5
1	B	1047	CYS	2.5
1	B	748	HIS	2.5
1	B	594	PHE	2.4
1	A	68	ASP	2.4
1	A	717	LYS	2.4
1	B	219	GLN	2.3
1	B	750	VAL	2.3
1	B	743	SER	2.3
1	B	593	LYS	2.3
1	A	298	LEU	2.3
1	B	691	PRO	2.2
1	B	676	THR	2.2
1	B	746	VAL	2.2
1	B	63	ARG	2.2
1	A	396	LEU	2.2
1	B	896	GLU	2.2
1	B	592	LEU	2.2
1	A	112	GLN	2.1
1	B	1012	LEU	2.1
1	A	305	VAL	2.1
1	B	539	ILE	2.1
1	A	1100	ALA	2.1
1	B	739	LEU	2.1
1	B	279	ILE	2.1
1	B	720	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	48	SER	2.1
1	B	763	CYS	2.1
1	A	718	LYS	2.0
1	A	747	TYR	2.0
1	B	893	PHE	2.0
1	A	470	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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	DOC	C	11	18/19	0.96	0.08	39,51,66,69	0
2	DOC	P	11	18/19	0.96	0.08	38,51,62,63	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

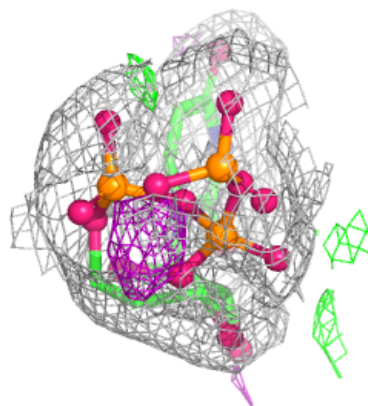
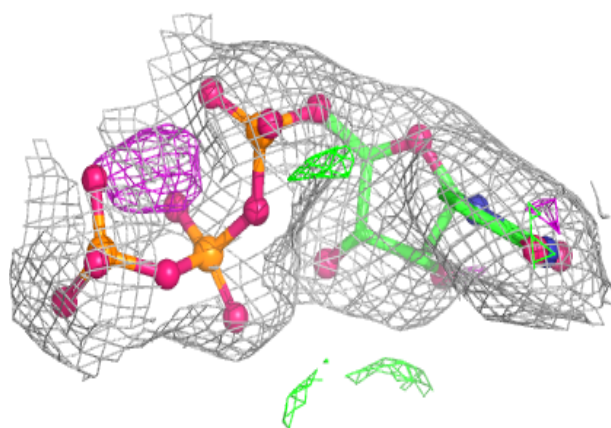
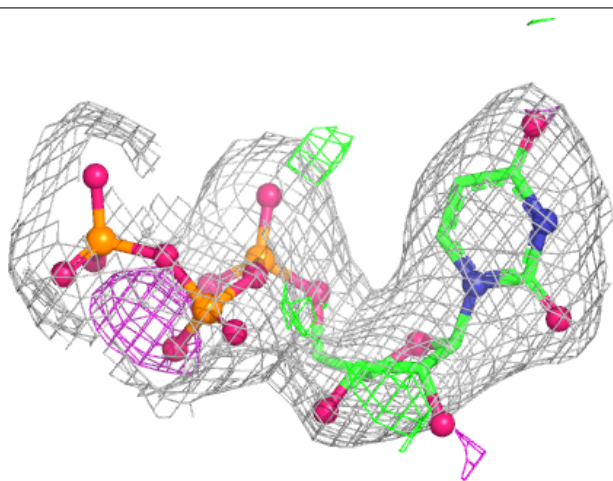
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

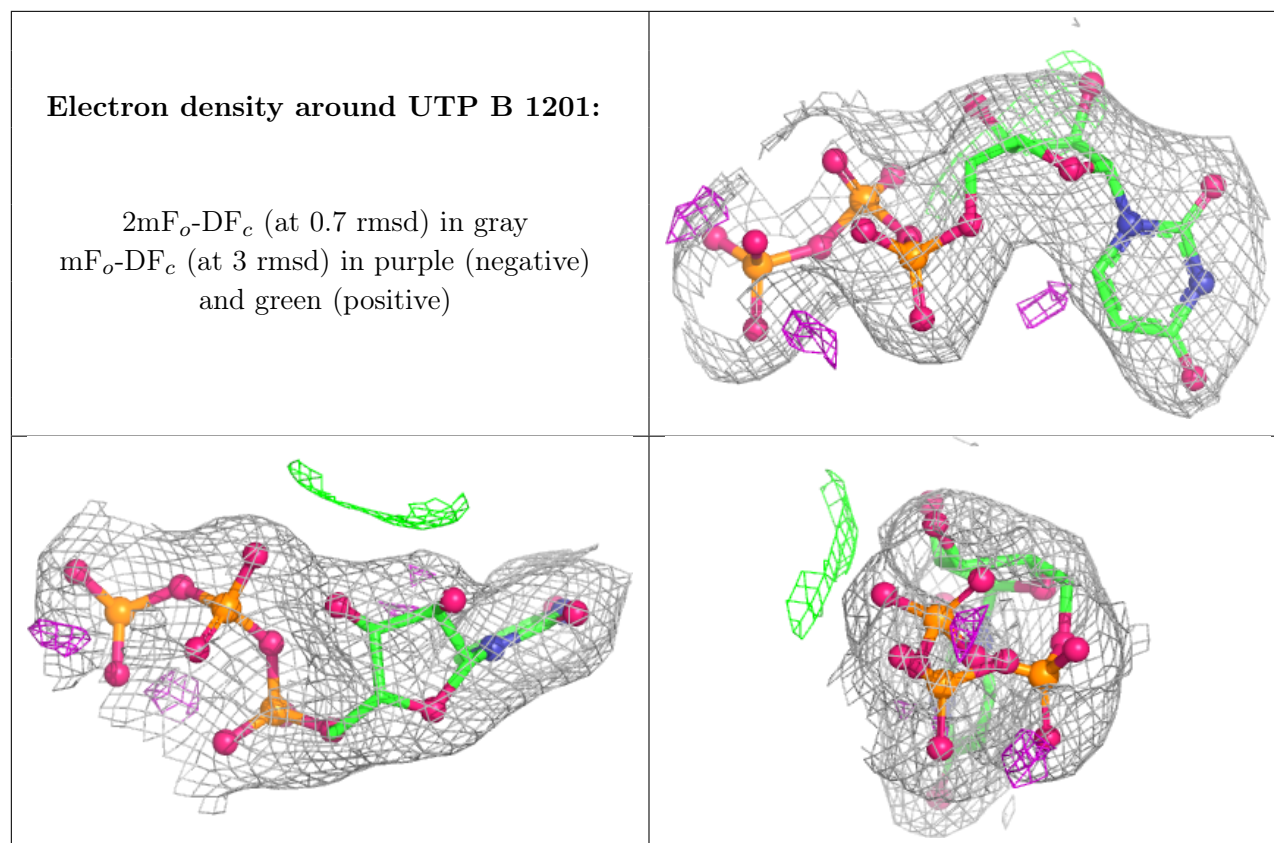
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	ACT	B	1205	4/4	0.83	0.18	72,74,77,77	0
6	ACT	A	1205	4/4	0.91	0.16	75,78,80,88	0
5	CA	B	1203	1/1	0.93	0.08	81,81,81,81	0
5	CA	A	1204	1/1	0.94	0.07	95,95,95,95	0
5	CA	A	1203	1/1	0.95	0.06	80,80,80,80	0
5	CA	A	1202	1/1	0.96	0.07	58,58,58,58	0
4	UTP	A	1201	29/29	0.96	0.08	44,53,60,64	0
4	UTP	B	1201	29/29	0.96	0.08	42,55,66,68	0
5	CA	B	1202	1/1	0.97	0.05	70,70,70,70	0
5	CA	B	1204	1/1	0.99	0.04	76,76,76,76	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 1201:

$2mF_o-DF_c$ (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.