



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:56 PM UTC

PDB ID : 3T64 / pdb_00003t64
Title : 5'-Diphenyl Nucleoside Inhibitors of Plasmodium falciparum dUTPase
Authors : Hampton, S.E.; Baragana, B.; Schipani, A.; Bosch-Navarrete, C.; Musso-Buendia, A.; Recio, E.; Kaiser, M.; Whittingham, J.L.; Roberts, S.M.; Shevtsov, M.; Brannigan, J.A.; Kahnberg, P.; Brun, R.; Wilson, K.S.; Gonzalez-Pacanowska, D.; Johansson, N.G.; Gilbert, I.H.
Deposited on : 2011-07-28
Resolution : 1.65 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](#) ①) 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

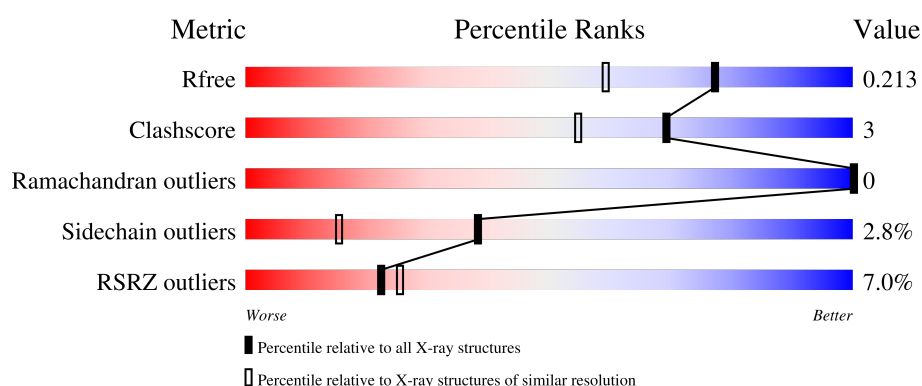
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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	2563 (1.66-1.66)
Clashscore	190562	2662 (1.66-1.66)
Ramachandran outliers	187476	2621 (1.66-1.66)
Sidechain outliers	187428	2621 (1.66-1.66)
RSRZ outliers	180081	2564 (1.66-1.66)

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	181	5% 74% 10% • 15%
1	B	181	5% 64% 10% • 25%
1	C	181	7% 74% 8% • 17%
2	F	3	33% 33% 67%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4043 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 Deoxyuridine 5'-triphosphate nucleotidohydrolase, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	153	Total	C	N	O	S	31	7	0
			1244	797	205	238	4			
1	B	135	Total	C	N	O	S	49	6	0
			1109	719	176	210	4			
1	C	151	Total	C	N	O	S	22	9	0
			1241	794	204	238	5			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	174	GLU	-	expression tag	UNP Q8II92
A	175	ALA	-	expression tag	UNP Q8II92
A	176	HIS	-	expression tag	UNP Q8II92
A	177	HIS	-	expression tag	UNP Q8II92
A	178	HIS	-	expression tag	UNP Q8II92
A	179	HIS	-	expression tag	UNP Q8II92
A	180	HIS	-	expression tag	UNP Q8II92
A	181	HIS	-	expression tag	UNP Q8II92
B	174	GLU	-	expression tag	UNP Q8II92
B	175	ALA	-	expression tag	UNP Q8II92
B	176	HIS	-	expression tag	UNP Q8II92
B	177	HIS	-	expression tag	UNP Q8II92
B	178	HIS	-	expression tag	UNP Q8II92
B	179	HIS	-	expression tag	UNP Q8II92
B	180	HIS	-	expression tag	UNP Q8II92
B	181	HIS	-	expression tag	UNP Q8II92
C	174	GLU	-	expression tag	UNP Q8II92
C	175	ALA	-	expression tag	UNP Q8II92
C	176	HIS	-	expression tag	UNP Q8II92
C	177	HIS	-	expression tag	UNP Q8II92
C	178	HIS	-	expression tag	UNP Q8II92
C	179	HIS	-	expression tag	UNP Q8II92
C	180	HIS	-	expression tag	UNP Q8II92

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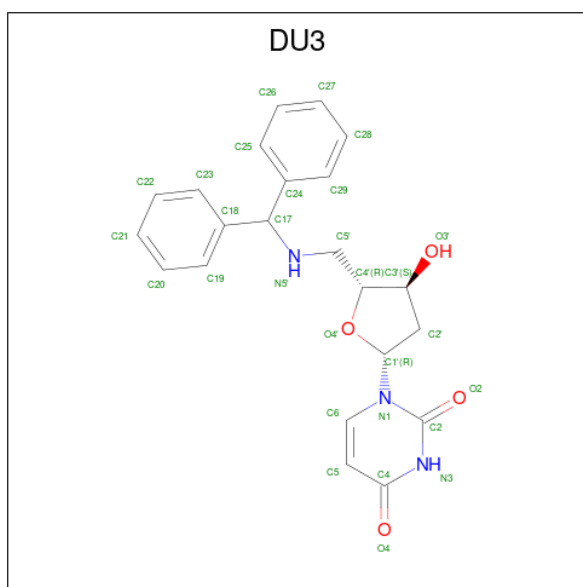
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Chain	Residue	Modelled	Actual	Comment	Reference
C	181	HIS	-	expression tag	UNP Q8II92

- Molecule 2 is a protein called 5'-(BENZHYDRYLAMINO)-2',5'-DIDEOXYURIDINE.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	F	3	Total	C	N	O	2	0	0
			20	12	5	3			

- Molecule 3 is 2',5'-dideoxy-5'-[(diphenylmethyl)amino]uridine (CCD ID: DU3) (formula: $C_{22}H_{23}N_3O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			29	22	3	4		
3	B	1	Total	C	N	O	0	0
			29	22	3	4		
3	C	1	Total	C	N	O	0	0
			29	22	3	4		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	134	Total	O	0	0
			134	134		
5	B	90	Total	O	0	0
			90	90		
5	C	96	Total	O	0	0
			96	96		
5	F	2	Total	O	0	0
			2	2		

- Molecule 1: Deoxyuridine 5'-triphosphate nucleotidohydrolase, putative



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	77.01Å 77.01Å 106.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.44 – 1.65 34.44 – 1.65	Depositor EDS
% Data completeness (in resolution range)	100.0 (34.44-1.65) 100.0 (34.44-1.65)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.88 (at 1.65Å)	Xtriage
Refinement program	REFMAC 5.5.0082	Depositor
R, R_{free}	0.190 , 0.216 0.188 , 0.213	Depositor DCC
R_{free} test set	3739 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	21.5	Xtriage
Anisotropy	0.206	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 42.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.038 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4043	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.23% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: DU3, SO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	1.01	7/1286 (0.5%)	0.89	5/1728 (0.3%)
1	B	0.95	5/1144 (0.4%)	0.83	0/1540
1	C	0.82	4/1284 (0.3%)	0.84	2/1724 (0.1%)
2	F	1.90	1/20 (5.0%)	2.60	2/26 (7.7%)
All	All	0.94	17/3734 (0.5%)	0.87	9/5018 (0.2%)

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	60[A]	SER	N-CA	-11.32	1.31	1.45
1	A	60[B]	SER	N-CA	-11.32	1.31	1.45
1	B	61	ASN	CA-CB	-8.45	1.41	1.53
2	F	1	ALA	CA-CB	-7.86	1.26	1.52
1	A	131	LYS	C-O	-6.21	1.16	1.24
1	A	132	LYS	C-O	-5.99	1.16	1.23
1	C	39	LEU	C-O	-5.85	1.16	1.24
1	A	126	GLN	C-O	-5.64	1.17	1.23
1	C	59	LYS	CB-CG	-5.46	1.36	1.52
1	A	130	ILE	C-O	-5.46	1.16	1.24
1	B	36	ASP	C-O	-5.39	1.17	1.23
1	B	12	VAL	C-O	-5.28	1.17	1.24
1	A	127	GLU	C-O	-5.25	1.17	1.23
1	B	132	LYS	CE-NZ	-5.19	1.33	1.49
1	B	27	SER	C-O	-5.18	1.17	1.24
1	C	15[A]	MET	C-O	-5.07	1.18	1.24
1	C	15[B]	MET	C-O	-5.07	1.18	1.24

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	24	GLU	CB-CA-C	9.07	123.45	109.84
2	F	1	ALA	N-CA-CB	8.75	123.52	110.40
2	F	3	ALA	CB-CA-C	-8.64	97.54	110.50
1	A	65	LYS	CG-CD-CE	-6.93	95.37	111.30
1	A	35	LYS	CG-CD-CE	5.82	124.67	111.30
1	C	20	LYS	CG-CD-CE	-5.72	98.15	111.30
1	C	131	LYS	CB-CG-CD	-5.69	98.22	111.30
1	A	18	ASN	N-CA-CB	-5.55	103.15	111.25
1	A	18	ASN	CB-CA-C	-5.38	102.09	111.23

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	1244	0	1276	6	1
1	B	1109	0	1154	10	0
1	C	1241	0	1274	9	1
2	F	20	0	19	0	0
3	A	29	0	23	1	0
3	B	29	0	23	0	0
3	C	29	0	23	0	0
4	A	5	0	0	0	0
4	B	5	0	0	0	0
4	C	10	0	0	0	0
5	A	134	0	0	1	1
5	B	90	0	0	0	0
5	C	96	0	0	1	1
5	F	2	0	0	0	0
All	All	4043	0	3792	25	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (25) 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:C:21:THR:HG21	1:C:30:ASP:OD2	1.60	1.01
1:B:1[A]:MET:HE2	1:B:147[A]:LEU:HD11	1.52	0.91
1:C:15[A]:MET:HE2	1:C:132:LYS:HE3	1.49	0.91
1:C:57:GLN:HB2	1:C:84[B]:THR:HG21	1.76	0.68
1:B:12:VAL:HA	1:B:15:MET:HE3	1.76	0.67
1:A:15:MET:HE3	1:A:133:ASN:ND2	2.11	0.66
1:C:21:THR:CG2	1:C:30:ASP:OD2	2.42	0.59
1:B:1[A]:MET:HE2	1:B:147[A]:LEU:CD1	2.31	0.56
1:A:13[B]:ARG:NH1	5:A:431:HOH:O	2.38	0.56
1:C:15[B]:MET:HE3	1:C:133:ASN:ND2	2.21	0.56
1:B:12:VAL:HA	1:B:15:MET:CE	2.36	0.55
1:A:56:LEU:HD23	1:A:83:ASN:HD22	1.72	0.55
1:C:19:HIS:O	1:C:20:LYS:HB2	2.07	0.54
1:A:15:MET:HE3	1:A:133:ASN:HD22	1.71	0.54
1:C:15[B]:MET:HE3	1:C:133:ASN:HD22	1.76	0.51
1:A:56:LEU:CD2	1:A:83:ASN:HD22	2.24	0.51
1:B:15:MET:HE1	1:B:34:VAL:HG12	1.94	0.49
1:A:1[A]:MET:HE2	1:A:147:LEU:HD22	1.94	0.48
3:A:201:DU3:H19	1:B:96:LYS:NZ	2.28	0.48
1:B:19:HIS:HE1	1:B:30:ASP:O	1.98	0.47
1:B:55:ALA:O	1:B:83:ASN:HA	2.16	0.46
1:C:132:LYS:O	1:C:133:ASN:HB2	2.17	0.44
1:B:84[B]:THR:HG22	5:C:387:HOH:O	2.18	0.44
1:C:1[B]:MET:HG3	1:C:57:GLN:HE21	1.83	0.44
1:B:15:MET:HE1	1:B:34:VAL:HB	2.02	0.41

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:96:LYS:NZ	5:C:388:HOH:O[4_564]	0.73	1.47
1:A:96:LYS:NZ	5:A:423:HOH:O[4_554]	2.14	0.06

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	155/181 (86%)	150 (97%)	5 (3%)	0	100	100
1	B	134/181 (74%)	131 (98%)	3 (2%)	0	100	100
1	C	153/181 (84%)	145 (95%)	8 (5%)	0	100	100
2	F	1/3 (33%)	1 (100%)	0	0	100	100
All	All	443/546 (81%)	427 (96%)	16 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	143/163 (88%)	141 (99%)	2 (1%)	59	39
1	B	129/163 (79%)	123 (95%)	6 (5%)	23	5
1	C	143/163 (88%)	138 (96%)	5 (4%)	32	10
2	F	1/1 (100%)	1 (100%)	0	100	100
All	All	416/490 (85%)	403 (97%)	13 (3%)	38	12

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	ASP
1	A	24	GLU
1	B	10	ASP
1	B	17	LYS
1	B	26	ASP
1	B	109[A]	ASP
1	B	109[B]	ASP
1	B	131	LYS
1	C	35	LYS
1	C	59	LYS

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Mol	Chain	Res	Type
1	C	61[A]	ASN
1	C	61[B]	ASN
1	C	109	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	HIS
1	A	83	ASN
1	A	133	ASN
1	B	19	HIS
1	C	22	HIS
1	C	57	GLN
1	C	126	GLN
1	C	133	ASN

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](#)

7 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	DU3	C	201	-	31,32,32	0.97	1 (3%)	39,44,44	1.63	6 (15%)
4	SO4	C	203	-	4,4,4	0.23	0	6,6,6	0.12	0
4	SO4	A	202	-	4,4,4	0.27	0	6,6,6	0.33	0
4	SO4	B	202	-	4,4,4	0.23	0	6,6,6	0.21	0
3	DU3	B	201	-	31,32,32	1.22	2 (6%)	39,44,44	1.54	4 (10%)
3	DU3	A	201	-	31,32,32	1.08	2 (6%)	39,44,44	1.78	7 (17%)
4	SO4	C	202	-	4,4,4	0.25	0	6,6,6	0.17	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
3	DU3	C	201	-	-	4/17/29/29	0/4/4/4
3	DU3	B	201	-	-	4/17/29/29	0/4/4/4
3	DU3	A	201	-	-	4/17/29/29	0/4/4/4

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	201	DU3	C24-C17	3.29	1.56	1.52
3	A	201	DU3	C5'-C4'	3.25	1.59	1.52
3	B	201	DU3	C18-C17	2.95	1.56	1.52
3	A	201	DU3	C6-C5	2.59	1.41	1.35
3	C	201	DU3	C6-C5	2.13	1.40	1.35

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	201	DU3	C4-N3-C2	-5.71	119.52	126.61
3	C	201	DU3	C4-N3-C2	-5.65	119.60	126.61
3	A	201	DU3	C4-N3-C2	-5.43	119.88	126.61
3	A	201	DU3	C5-C4-N3	4.65	121.32	114.80
3	C	201	DU3	C5-C4-N3	4.36	120.91	114.80
3	B	201	DU3	C5-C4-N3	4.00	120.40	114.80
3	A	201	DU3	C24-C17-C18	-3.83	104.64	112.33
3	B	201	DU3	N3-C2-N1	3.61	119.60	114.89
3	A	201	DU3	O4'-C1'-N1	-3.56	101.54	107.86
3	A	201	DU3	N3-C2-N1	3.42	119.35	114.89
3	C	201	DU3	N3-C2-N1	3.13	118.97	114.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	201	DU3	O4'-C1'-N1	-3.07	102.41	107.86
3	B	201	DU3	O4-C4-C5	-2.46	120.91	125.16
3	C	201	DU3	C25-C24-C17	-2.45	116.62	120.75
3	A	201	DU3	O4-C4-C5	-2.27	121.25	125.16
3	C	201	DU3	O4-C4-C5	-2.22	121.33	125.16
3	A	201	DU3	C5'-C4'-C3'	-2.14	111.06	115.55

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	201	DU3	C2'-C1'-N1-C6
3	B	201	DU3	C2'-C1'-N1-C6
3	C	201	DU3	C2'-C1'-N1-C6
3	A	201	DU3	C2'-C1'-N1-C2
3	C	201	DU3	C2'-C1'-N1-C2
3	A	201	DU3	O4'-C1'-N1-C6
3	B	201	DU3	O4'-C1'-N1-C6
3	C	201	DU3	O4'-C1'-N1-C6
3	B	201	DU3	C2'-C1'-N1-C2
3	A	201	DU3	O4'-C1'-N1-C2
3	B	201	DU3	O4'-C1'-N1-C2
3	C	201	DU3	O4'-C1'-N1-C2

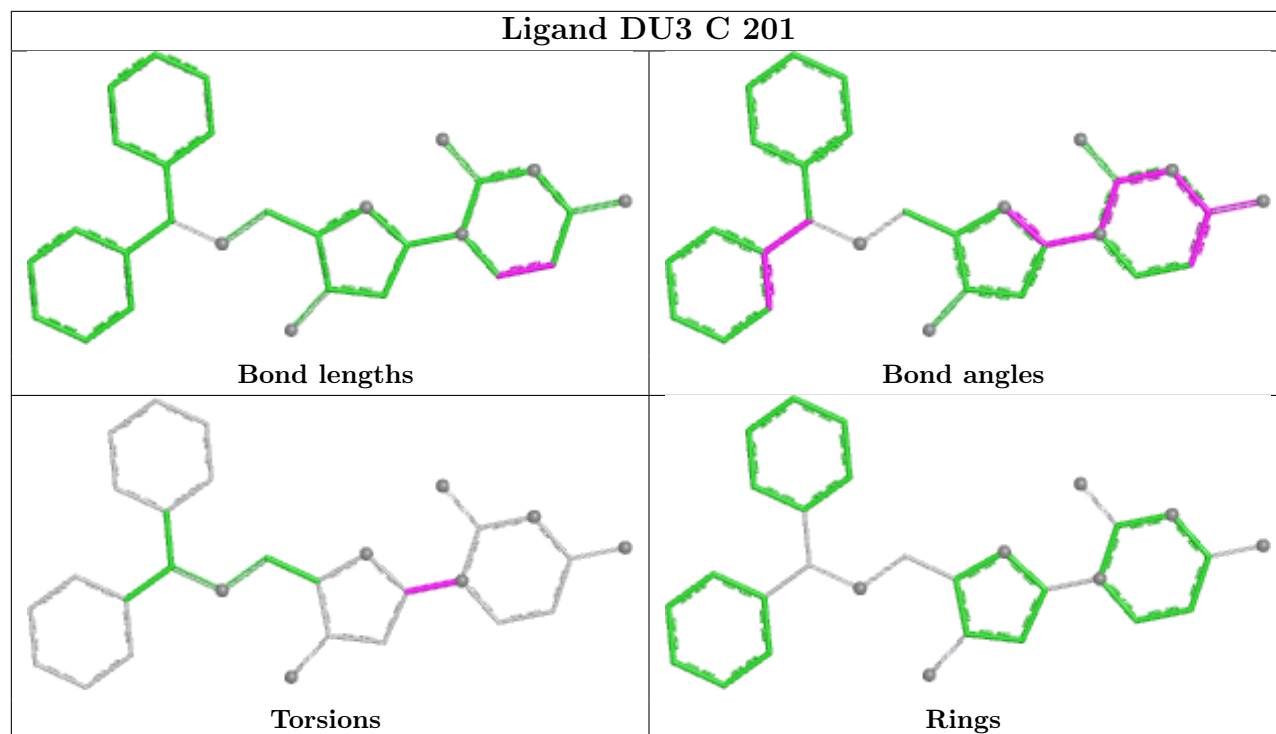
There are no ring outliers.

1 monomer is involved in 1 short contact:

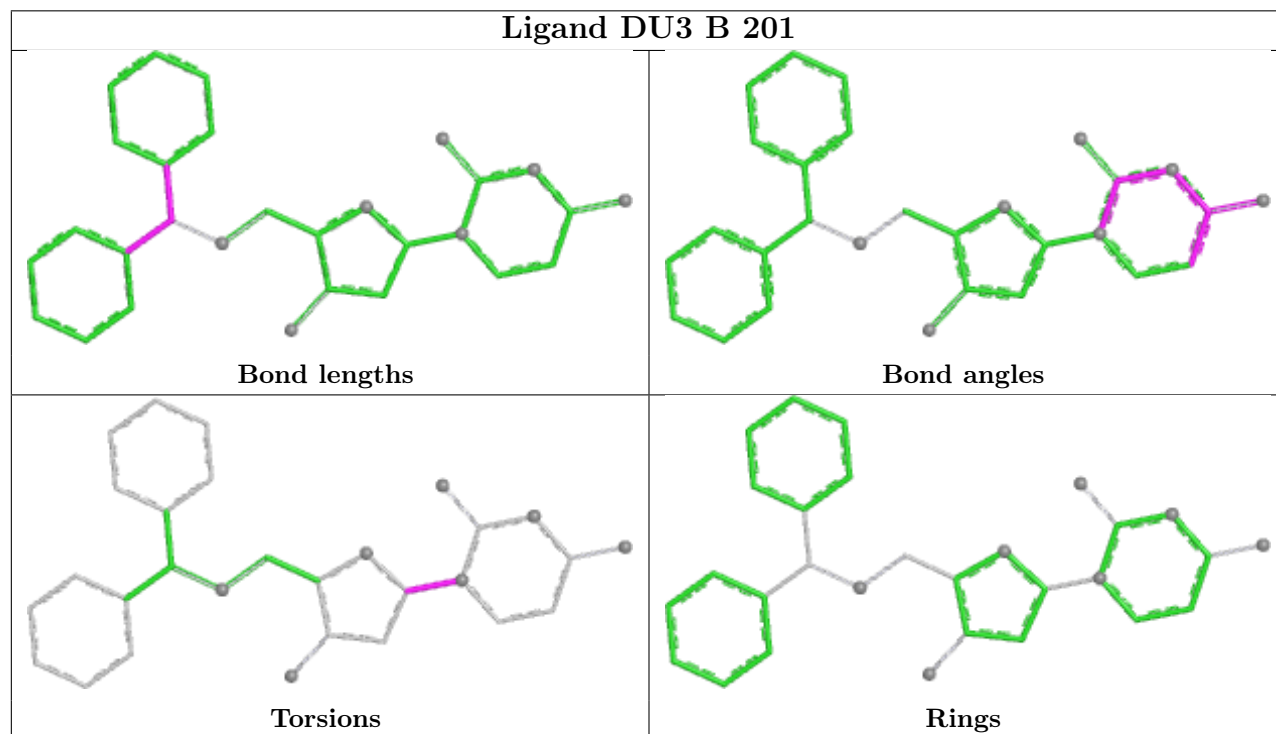
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	201	DU3	1	0

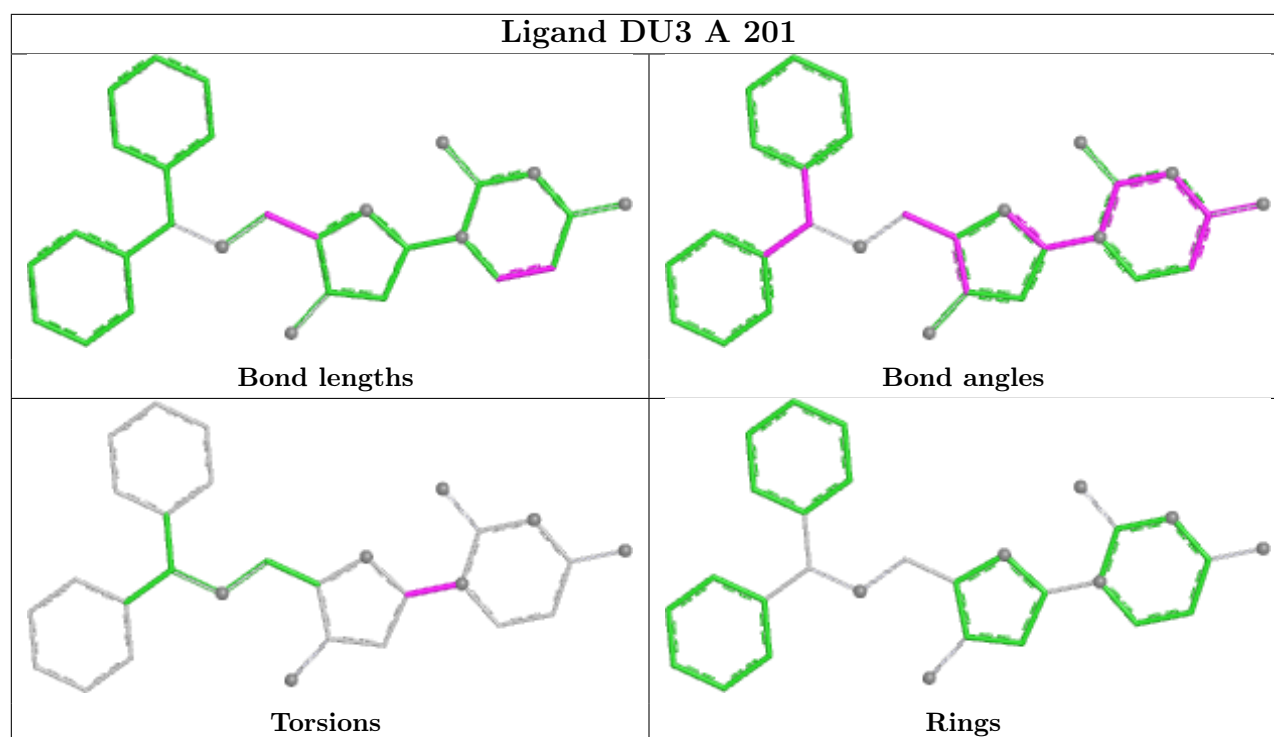
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 DU3 C 201



Ligand DU3 B 201





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	153/181 (84%)	-0.01	9 (5%)	28 32	8, 18, 32, 44	18 (11%)
1	B	135/181 (74%)	0.30	9 (6%)	24 27	11, 23, 38, 43	19 (14%)
1	C	151/181 (83%)	0.23	12 (7%)	18 21	10, 20, 35, 54	17 (11%)
2	F	3/3 (100%)	1.69	1 (33%)	1 1	25, 25, 26, 30	2 (66%)
All	All	442/546 (80%)	0.18	31 (7%)	22 25	8, 20, 35, 54	56 (12%)

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	167	SER	5.6
1	C	21	THR	5.0
1	C	22	HIS	4.4
1	C	166	GLY	4.4
1	A	167	SER	3.6
1	B	18	ASN	3.1
1	C	81	ILE	3.0
1	A	23	HIS	2.9
1	C	20	LYS	2.9
1	A	80	ASN	2.9
1	C	19	HIS	2.9
1	A	25	GLY	2.7
1	C	25	GLY	2.7
1	B	64	TYR	2.6
1	A	81	ILE	2.6
1	C	26	ASP	2.6
1	A	83	ASN	2.5
1	B	155	LEU	2.4
1	A	22	HIS	2.4
1	C	18[A]	ASN	2.3
1	C	64	TYR	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	10	ASP	2.3
1	B	151	LEU	2.3
1	A	65	LYS	2.2
1	A	64	TYR	2.2
1	B	19	HIS	2.2
1	B	65	LYS	2.2
1	B	81	ILE	2.1
1	C	10	ASP	2.1
1	B	62	TYR	2.0
2	F	3	ALA	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](#)

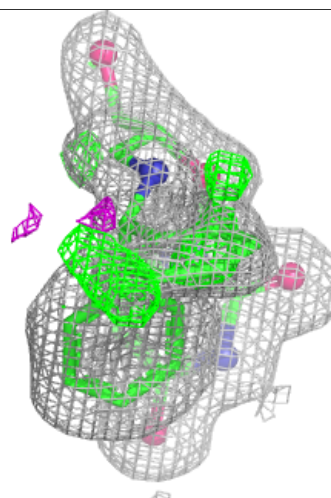
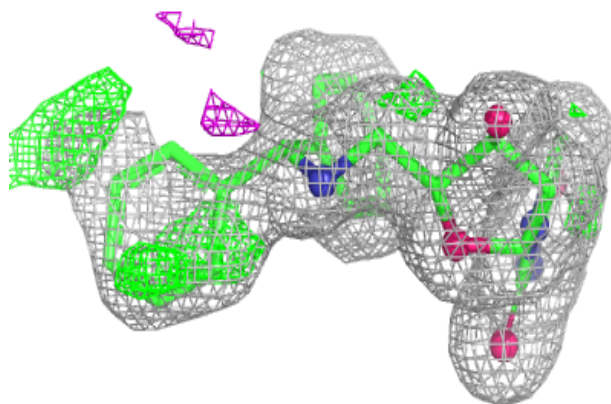
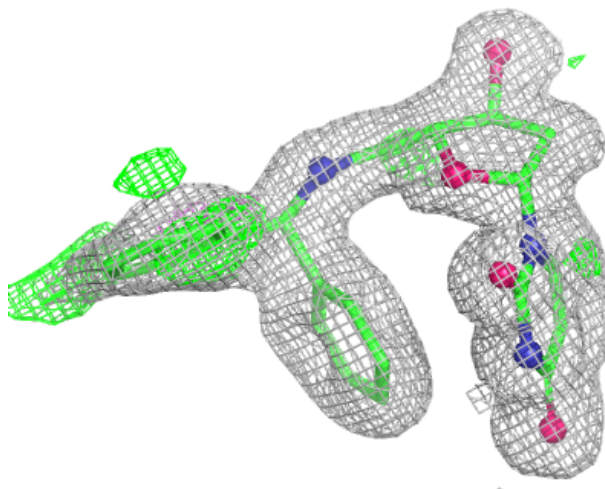
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
4	SO4	C	203	5/5	0.83	0.12	55,55,55,56	0
4	SO4	C	202	5/5	0.89	0.12	46,47,47,48	0
4	SO4	B	202	5/5	0.91	0.10	37,38,40,40	0
4	SO4	A	202	5/5	0.92	0.10	38,39,40,40	0
3	DU3	A	201	29/29	0.95	0.10	14,23,30,30	6
3	DU3	B	201	29/29	0.95	0.09	16,24,31,31	6
3	DU3	C	201	29/29	0.96	0.10	13,23,29,29	6

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.

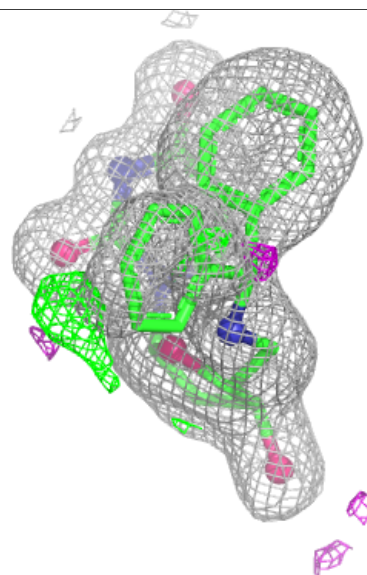
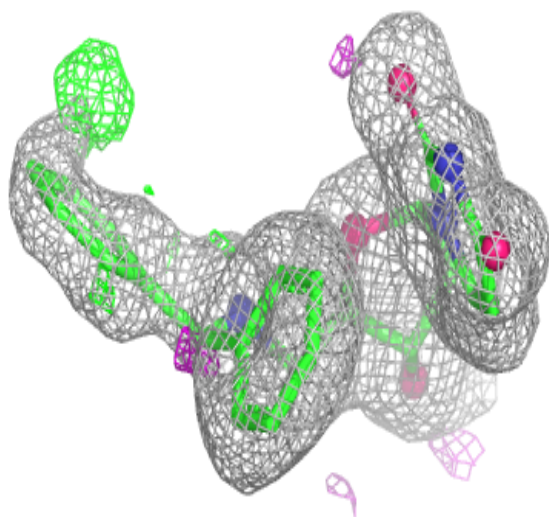
Electron density around DU3 A 201:

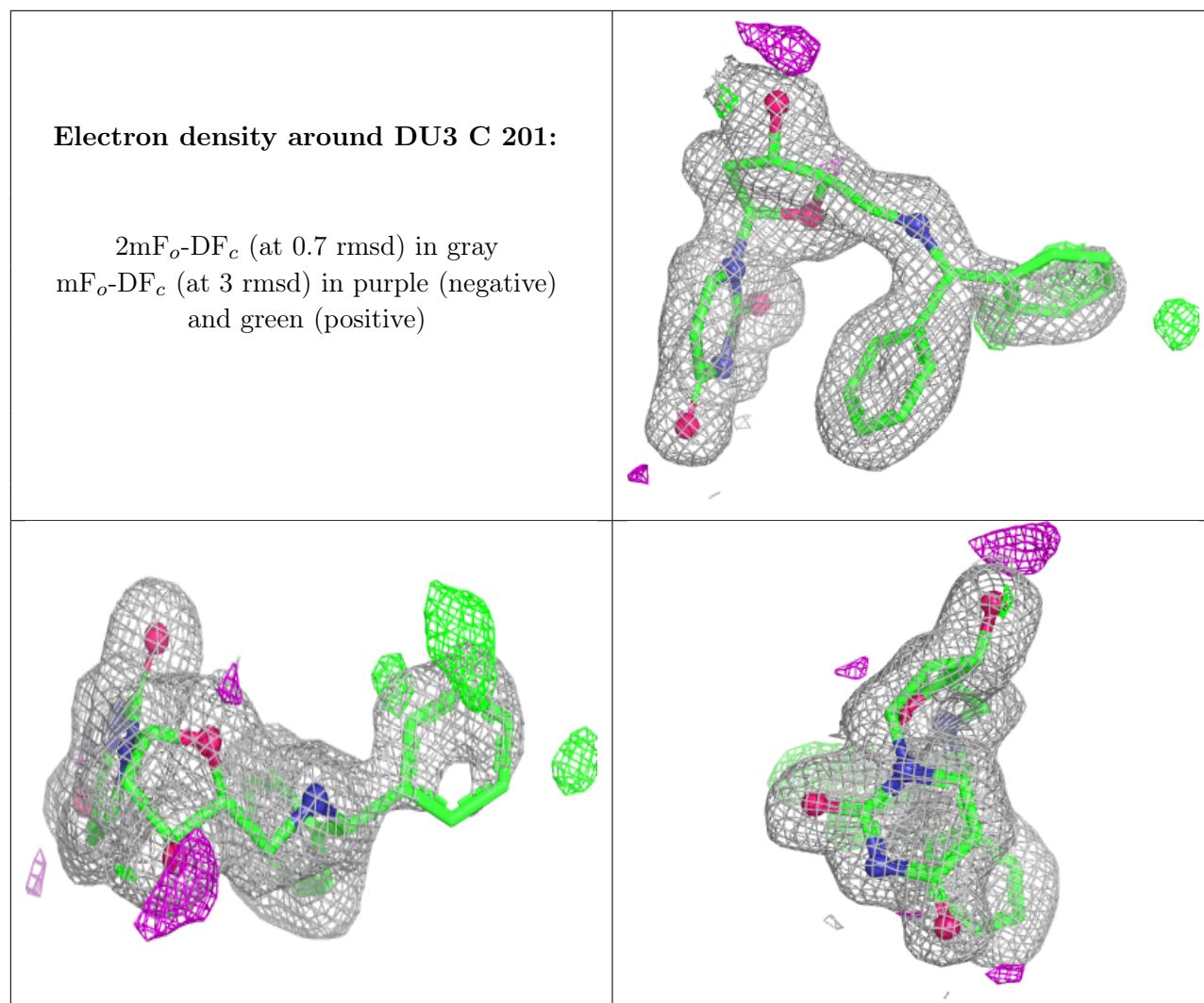
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around DU3 B 201:

$2mF_o-DF_c$ (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.