



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 06:36 PM UTC

PDB ID : 3TI0 / pdb_00003ti0
Title : Crystal Structure of Bacillus DNA Polymerase I Large Fragment Bound to DNA and ddGTP-dC in Closed Conformation
Authors : Wang, W.; Beese, L.S.
Deposited on : 2011-08-19
Resolution : 1.62 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

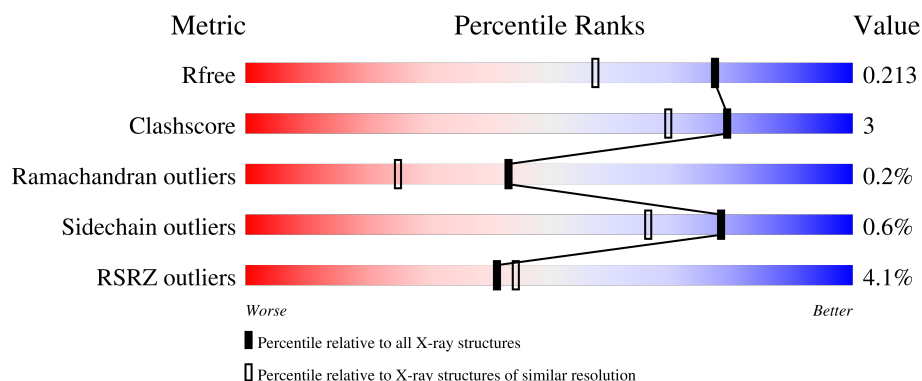
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.62 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	6728 (1.64-1.60)
Clashscore	190562	7023 (1.64-1.60)
Ramachandran outliers	187476	6898 (1.64-1.60)
Sidechain outliers	187428	6896 (1.64-1.60)
RSRZ outliers	180081	6727 (1.64-1.60)

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	592	7% 91% 7%
1	D	592	% 91% 6%
2	B	9	78% 11% 11%
2	E	9	89% 11%
3	C	13	8% 69% 23% 8%

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Mol	Chain	Length	Quality of chain
3	F	13	8% 92% 8%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 21902 atoms, of which 9966 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 I.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	579	Total	C	H	N	O	S	0	2	0
			9378	2959	4724	809	869	17			
1	D	579	Total	C	H	N	O	S	0	10	0
			9427	2971	4748	813	878	17			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	456	GLU	ALA	SEE REMARK 999	UNP C9RTX7
A	505	LYS	GLU	SEE REMARK 999	UNP C9RTX7
A	598	ALA	ASP	engineered mutation	UNP C9RTX7
A	710	TYR	PHE	engineered mutation	UNP C9RTX7
A	823	HIS	ARG	SEE REMARK 999	UNP C9RTX7
D	456	GLU	ALA	SEE REMARK 999	UNP C9RTX7
D	505	LYS	GLU	SEE REMARK 999	UNP C9RTX7
D	598	ALA	ASP	engineered mutation	UNP C9RTX7
D	710	TYR	PHE	engineered mutation	UNP C9RTX7
D	823	HIS	ARG	SEE REMARK 999	UNP C9RTX7

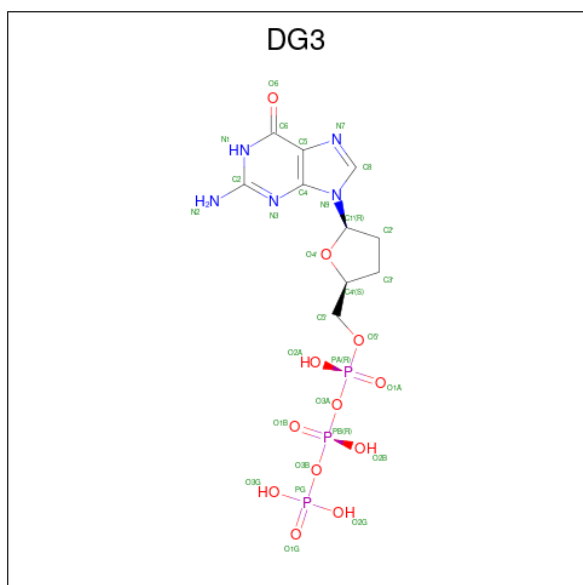
- Molecule 2 is a DNA chain called 5'-D(*CP*CP*TP*GP*AP*CP*TP*CP*(DDG))-3'.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	9	Total	C	H	N	O	P	0	0	0
			278	86	101	31	52	8			
2	E	9	Total	C	H	N	O	P	0	0	0
			278	86	101	31	52	8			

- Molecule 3 is a DNA chain called 5'-D(*C*AP*TP*CP*CP*GP*AP*GP*TP*CP*AP*GP*G)-3'.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	C	12	Total	C	H	N	O	P	0	0	0
			380	117	135	48	69	11			
3	F	12	Total	C	H	N	O	P	0	0	0
			380	117	135	48	69	11			

- Molecule 4 is 2'-3'-DIDEOXYGUANOSINE-5'-TRIPHOSPHATE (CCD ID: DG3) (formula: $C_{10}H_{16}N_5O_{12}P_3$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	H	N	O	P	0	0
			41	10	11	5	12	3		
4	D	1	Total	C	H	N	O	P	0	0
			41	10	11	5	12	3		

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Mg	0	0
			1	1		
5	D	1	Total	Mg	0	0
			1	1		

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	S	0	0
			5	4	1		
6	D	1	Total	O	S	0	0
			5	4	1		

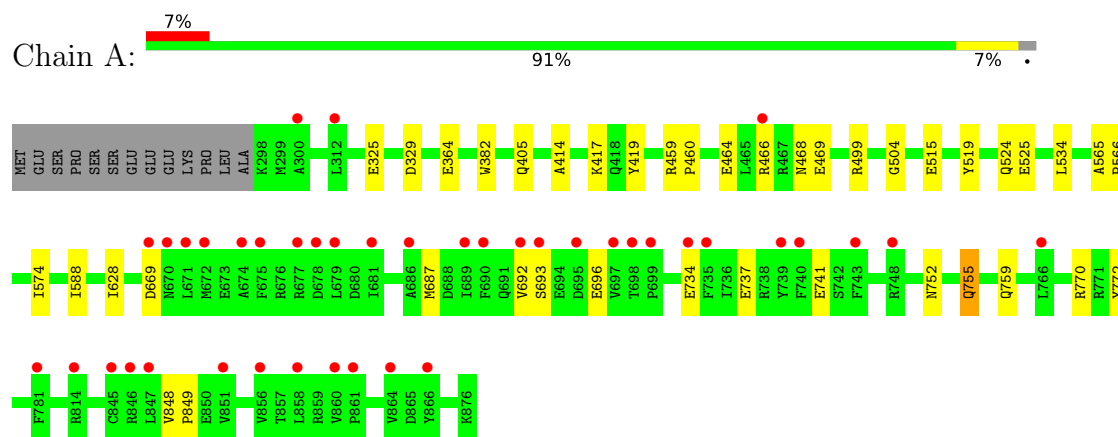
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	587	Total	O	0	0
			587	587		
7	D	854	Total	O	0	0
			854	854		
7	B	38	Total	O	0	0
			38	38		
7	C	78	Total	O	0	0
			78	78		
7	E	40	Total	O	0	0
			40	40		
7	F	90	Total	O	0	0
			90	90		

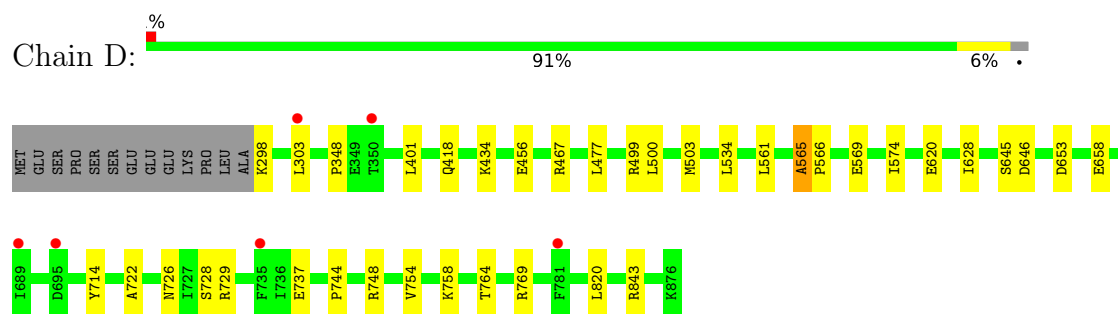
3 Residue-property plots [i](#)

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.

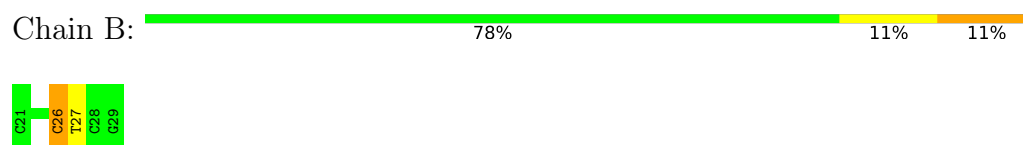
• Molecule 1: DNA polymerase I



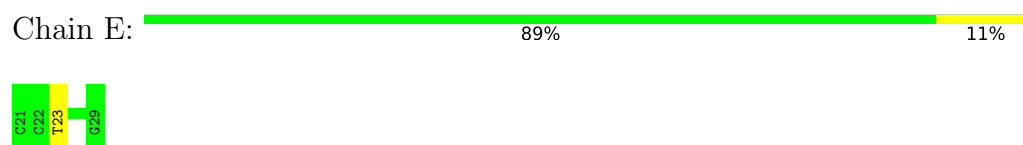
• Molecule 1: DNA polymerase I



• Molecule 2: 5'-D(*CP*CP*TP*GP*AP*CP*TP*CP*(DDG))-3'

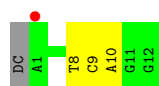


• Molecule 2: 5'-D(*CP*CP*TP*GP*AP*CP*TP*CP*(DDG))-3'



- Molecule 3: 5'-D(*C*AP*TP*CP*CP*GP*AP*GP*TP*CP*AP*GP*G)-3'

Chain C:  8% 69% 23% 8%



- Molecule 3: 5'-D(*C*AP*TP*CP*CP*GP*AP*GP*TP*CP*AP*GP*G)-3'

Chain F:  8% 92% 8%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	93.50Å 108.92Å 149.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.96 – 1.62 42.96 – 1.62	Depositor EDS
% Data completeness (in resolution range)	93.4 (42.96-1.62) 93.4 (42.96-1.62)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.61 (at 1.62Å)	Xtriage
Refinement program	PHENIX 1.7.1 _743	Depositor
R, R_{free}	0.187 , 0.215 0.185 , 0.213	Depositor DCC
R_{free} test set	8205 reflections (4.54%)	wwPDB-VP
Wilson B-factor (Å ²)	17.5	Xtriage
Anisotropy	0.112	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 43.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	21902	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.99% 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: DG3, MG, DDG, 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	0.44	0/4747	0.70	0/6416
1	D	0.56	0/4802	0.77	3/6488 (0.0%)
2	B	0.41	0/173	1.23	1/264 (0.4%)
2	E	0.44	0/173	1.19	1/264 (0.4%)
3	C	0.39	0/275	0.98	1/423 (0.2%)
3	F	0.43	0/275	1.02	0/423
All	All	0.50	0/10445	0.78	6/14278 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	26	DC	N1-C1'-C2'	-7.39	102.41	113.50
3	C	8	DT	O4'-C1'-N1	-5.48	100.18	108.40
1	D	620	GLU	CB-CA-C	-5.44	106.27	111.00
2	E	23	DT	O4'-C1'-N1	-5.20	100.60	108.40
1	D	565	ALA	CA-C-N	-5.17	114.50	119.87
1	D	565	ALA	C-N-CA	-5.17	114.50	119.87

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	4654	4724	4703	27	0
1	D	4679	4748	4694	26	0
2	B	177	101	103	1	0
2	E	177	101	103	0	0
3	C	245	135	136	1	0
3	F	245	135	136	0	0
4	A	30	11	12	0	0
4	D	30	11	12	1	0
5	A	1	0	0	0	0
5	D	1	0	0	0	0
6	A	5	0	0	0	0
6	D	5	0	0	0	0
7	A	587	0	0	7	0
7	B	38	0	0	0	0
7	C	78	0	0	0	0
7	D	854	0	0	15	0
7	E	40	0	0	0	0
7	F	90	0	0	0	0
All	All	11936	9966	9899	55	0

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

All (55) 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:752:ASN:O	1:A:755:GLN:HG3	1.92	0.69
1:A:499:ARG:NE	7:A:1008:HOH:O	2.28	0.66
1:A:405:GLN:NE2	7:A:1672:HOH:O	2.31	0.63
1:A:848:VAL:HB	1:A:849:PRO:HD3	1.81	0.61
1:D:434:LYS:NZ	7:D:1040:HOH:O	2.34	0.60
1:D:569:GLU:CD	7:D:1107:HOH:O	2.45	0.59
1:A:466:ARG:HD2	7:A:1627:HOH:O	2.02	0.59
3:C:9:DC:H2''	3:C:10:DA:C8	2.38	0.58
1:A:459:ARG:HB3	1:A:460:PRO:HD3	1.86	0.58
1:D:653:ASP:OD2	7:D:1686:HOH:O	2.17	0.58
1:D:748:ARG:HD3	7:D:1129:HOH:O	2.06	0.55
1:D:499:ARG:NH2	7:D:1527:HOH:O	2.39	0.55
1:D:418:GLN:HA	7:D:1059:HOH:O	2.07	0.54
1:D:754:VAL:CG1	1:D:758:LYS:HE2	2.38	0.53
1:A:770:ARG:HD3	1:A:772:TYR:CZ	2.43	0.53
1:D:754:VAL:HG12	1:D:758:LYS:HE2	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:348:PRO:HB2	7:D:1161:HOH:O	2.09	0.52
1:A:692:VAL:O	1:A:693:SER:OG	2.22	0.51
1:A:325:GLU:H	1:A:325:GLU:CD	2.19	0.50
1:D:645:SER:O	1:D:646:ASP:HB2	2.13	0.49
1:A:687:MET:HG2	1:A:693:SER:C	2.38	0.48
1:D:744:PRO:O	1:D:748:ARG:HG3	2.13	0.48
1:A:459:ARG:NH2	7:A:1550:HOH:O	2.46	0.47
1:D:500:LEU:HA	1:D:503:MET:HE3	1.95	0.47
1:D:714:TYR:HB3	7:D:1032:HOH:O	2.14	0.47
1:A:692:VAL:HG22	1:A:696:GLU:HB2	1.96	0.47
1:D:298:LYS:HB2	7:D:1586:HOH:O	2.14	0.47
1:A:414:ALA:HB1	1:A:419:TYR:HB3	1.97	0.46
1:D:658:GLU:CD	4:D:202:DG3:H2'1	2.40	0.46
1:D:456:GLU:HG3	7:D:1175:HOH:O	2.16	0.46
1:D:534:LEU:HD11	1:D:574:ILE:HD13	1.97	0.46
1:A:499:ARG:CZ	7:A:1008:HOH:O	2.64	0.45
1:D:565:ALA:N	1:D:566:PRO:CD	2.80	0.45
1:A:565:ALA:N	1:A:566:PRO:CD	2.81	0.44
1:A:759:GLN:HG2	7:D:1465:HOH:O	2.18	0.43
1:A:534:LEU:HD11	1:A:574:ILE:HD13	2.01	0.43
1:A:504:GLY:HA2	1:A:588:ILE:HD12	2.01	0.43
1:D:653:ASP:HB2	7:D:1687:HOH:O	2.18	0.43
1:A:417:LYS:NZ	1:A:464:GLU:OE2	2.40	0.42
1:A:329:ASP:H	1:A:382:TRP:CD1	2.37	0.42
2:B:26:DC:H2''	2:B:27:DT:O5'	2.19	0.42
1:D:737:GLU:HG3	7:D:1493:HOH:O	2.19	0.42
1:A:524:GLN:HG2	1:A:525:GLU:O	2.20	0.42
1:A:364:GLU:HG3	7:A:207:HOH:O	2.20	0.41
1:A:515:GLU:HG2	1:A:519:TYR:CE2	2.55	0.41
1:D:467:ARG:NH2	7:D:1626:HOH:O	2.27	0.41
1:D:764:THR:HA	1:D:769:ARG:O	2.20	0.41
1:D:820:LEU:HD21	1:D:843:ARG:CZ	2.50	0.41
1:A:468:ASN:O	1:A:469:GLU:HB2	2.21	0.41
1:D:722:ALA:HB2	1:D:729:ARG:HA	2.03	0.41
1:D:726:ASN:HB2	7:D:1121:HOH:O	2.21	0.41
1:D:401:LEU:HD23	1:D:401:LEU:HA	1.93	0.40
1:A:499:ARG:CG	7:A:1008:HOH:O	2.68	0.40
1:A:737:GLU:O	1:A:741:GLU:HG3	2.21	0.40
1:A:734:GLU:HA	1:A:734:GLU:OE1	2.21	0.40

There are no symmetry-related clashes.

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	579/592 (98%)	561 (97%)	17 (3%)	1 (0%)	43	25
1	D	587/592 (99%)	574 (98%)	12 (2%)	1 (0%)	43	25
All	All	1166/1184 (98%)	1135 (97%)	29 (2%)	2 (0%)	43	25

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	628	ILE
1	A	628	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	497/507 (98%)	495 (100%)	2 (0%)	84	75
1	D	505/507 (100%)	501 (99%)	4 (1%)	73	59
All	All	1002/1014 (99%)	996 (99%)	6 (1%)	78	66

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	669	ASP
1	A	755	GLN
1	D	303	LEU
1	D	477	LEU

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Mol	Chain	Res	Type
1	D	561	LEU
1	D	728	SER

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

Mol	Chain	Res	Type
1	D	418	GLN
1	D	700	ASN
1	D	704	GLN

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	DDG	B	29	2,3	20,23,24	0.73	0	27,33,36	0.68	0
2	DDG	E	29	2,3	20,23,24	0.53	0	27,33,36	0.52	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	DDG	B	29	2,3	-	0/7/18/19	0/3/3/3
2	DDG	E	29	2,3	-	0/7/18/19	0/3/3/3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 2 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	$\# Z > 2$
4	DG3	A	201	5	31,32,32	2.12	6 (19%)	44,50,50	2.04	15 (34%)
6	SO4	A	2	-	4,4,4	0.27	0	6,6,6	0.12	0
6	SO4	D	1	-	4,4,4	0.23	0	6,6,6	0.24	0
4	DG3	D	202	5	31,32,32	1.86	6 (19%)	44,50,50	1.78	13 (29%)

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	DG3	A	201	5	-	6/22/31/31	0/3/3/3
4	DG3	D	202	5	-	5/22/31/31	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	201	DG3	O6-C6	7.24	1.37	1.23
4	A	201	DG3	C2-N2	5.91	1.48	1.34
4	D	202	DG3	O6-C6	5.34	1.33	1.23
4	D	202	DG3	C2-N2	4.98	1.45	1.34
4	D	202	DG3	PB-O3A	2.90	1.62	1.59
4	D	202	DG3	C1'-N9	-2.82	1.41	1.48
4	A	201	DG3	C1'-N9	-2.68	1.41	1.48
4	A	201	DG3	C2'-C3'	-2.68	1.46	1.54
4	D	202	DG3	C6-N1	-2.42	1.34	1.38
4	A	201	DG3	PB-O3A	2.35	1.62	1.59
4	D	202	DG3	C2'-C3'	-2.20	1.48	1.54
4	A	201	DG3	C6-N1	-2.11	1.34	1.38

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	201	DG3	C5-C4-N3	-4.96	120.49	128.39
4	A	201	DG3	O4'-C1'-N9	4.57	115.97	107.86
4	A	201	DG3	C2-N3-C4	4.19	119.52	112.30
4	D	202	DG3	C5-C4-N3	-3.77	122.39	128.39
4	A	201	DG3	C2'-C1'-N9	-3.63	105.52	112.40
4	A	201	DG3	O2B-PB-O3B	3.56	116.91	107.27
4	D	202	DG3	C2'-C1'-N9	-3.56	105.65	112.40
4	D	202	DG3	C6-C5-N7	3.35	136.38	130.29
4	A	201	DG3	C6-C5-N7	3.26	136.22	130.29
4	D	202	DG3	C2-N1-C6	-3.23	119.26	125.11
4	A	201	DG3	O3A-PB-O1B	-3.20	101.09	110.70
4	D	202	DG3	O4'-C1'-N9	3.16	113.47	107.86
4	D	202	DG3	C2-N3-C4	3.06	117.56	112.30
4	A	201	DG3	O2G-PG-O3B	2.93	114.45	104.64
4	D	202	DG3	C5-C6-N1	2.81	120.40	113.25
4	A	201	DG3	C4-C5-N7	-2.69	106.41	110.67
4	A	201	DG3	C2-N1-C6	-2.44	120.69	125.11
4	A	201	DG3	O2A-PA-O3A	2.42	113.82	107.27
4	A	201	DG3	C5-C6-N1	2.41	119.40	113.25
4	A	201	DG3	N9-C4-N3	2.39	130.73	125.95
4	D	202	DG3	O2B-PB-O3B	2.24	113.34	107.27
4	A	201	DG3	O2B-PB-O3A	2.22	113.29	107.27
4	D	202	DG3	O2G-PG-O3B	2.19	111.97	104.64
4	A	201	DG3	N9-C8-N7	-2.11	109.48	113.40
4	D	202	DG3	N9-C8-N7	-2.06	109.58	113.40
4	D	202	DG3	O6-C6-C5	-2.05	121.13	126.53
4	D	202	DG3	O2B-PB-O3A	2.04	112.78	107.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	202	DG3	C4-C5-N7	-2.01	107.49	110.67

There are no chirality outliers.

All (11) torsion outliers are listed below:

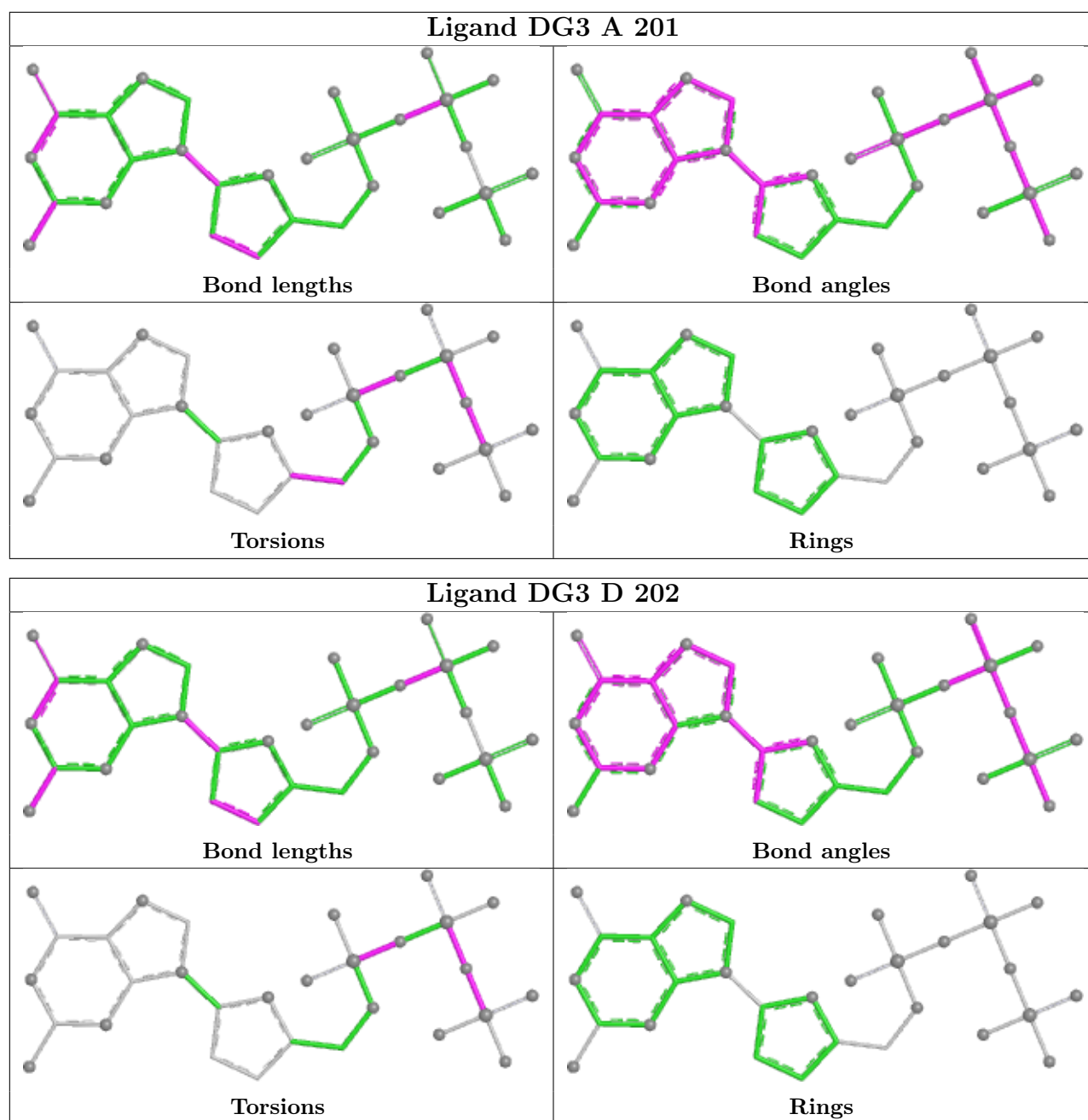
Mol	Chain	Res	Type	Atoms
4	D	202	DG3	PB-O3B-PG-O1G
4	A	201	DG3	PB-O3A-PA-O2A
4	D	202	DG3	PG-O3B-PB-O2B
4	D	202	DG3	PB-O3A-PA-O2A
4	A	201	DG3	PG-O3B-PB-O2B
4	A	201	DG3	PG-O3B-PB-O1B
4	A	201	DG3	O4'-C4'-C5'-O5'
4	A	201	DG3	PB-O3A-PA-O1A
4	D	202	DG3	PB-O3A-PA-O1A
4	A	201	DG3	PB-O3B-PG-O1G
4	D	202	DG3	PG-O3B-PB-O1B

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	202	DG3	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	579/592 (97%)	0.52	41 (7%) 22 23	9, 30, 53, 66	1 (0%)
1	D	579/592 (97%)	-0.18	6 (1%) 79 83	6, 19, 37, 49	5 (0%)
2	B	8/9 (88%)	-0.46	0 100 100	17, 18, 33, 36	0
2	E	8/9 (88%)	-0.42	0 100 100	12, 19, 31, 37	0
3	C	12/13 (92%)	-0.28	1 (8%) 17 18	15, 21, 41, 64	0
3	F	12/13 (92%)	-0.29	1 (8%) 17 18	11, 17, 36, 58	0
All	All	1198/1228 (97%)	0.15	49 (4%) 41 44	6, 24, 49, 66	6 (0%)

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	692	VAL	4.1
1	A	858	LEU	3.8
1	A	681	ILE	3.8
1	A	677	ARG	3.6
1	A	679	LEU	3.6
1	A	697	VAL	3.4
1	A	735	PHE	3.4
1	A	669	ASP	3.3
1	A	675	PHE	3.1
1	A	689	ILE	3.0
1	A	739	TYR	2.9
1	A	866	TYR	2.9
1	A	847	LEU	2.9
1	D	303	LEU	2.9
1	D	689	ILE	2.9
1	A	699	PRO	2.8
1	A	466	ARG	2.7
1	A	861	PRO	2.7
1	A	312	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	671	LEU	2.7
1	A	851	VAL	2.6
1	A	856	VAL	2.6
1	A	695	ASP	2.5
3	F	1	DA	2.5
1	A	300	ALA	2.4
1	A	693	SER	2.4
1	A	698	THR	2.4
1	A	860	VAL	2.4
1	A	686	ALA	2.3
1	A	674	ALA	2.3
1	A	690	PHE	2.3
1	A	678	ASP	2.3
1	A	670	ASN	2.3
1	A	814	ARG	2.3
3	C	1	DA	2.2
1	D	735	PHE	2.2
1	A	748	ARG	2.2
1	A	864	VAL	2.2
1	D	695	ASP	2.2
1	A	734	GLU	2.1
1	A	743	PHE	2.1
1	A	781	PHE	2.1
1	A	845	CYS	2.1
1	D	350	THR	2.1
1	A	740	PHE	2.1
1	A	766	LEU	2.0
1	A	672	MET	2.0
1	A	846	ARG	2.0
1	D	781	PHE	2.0

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

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	DDG	B	29	21/22	0.97	0.06	12,18,22,26	0
2	DDG	E	29	21/22	0.99	0.04	7,10,14,15	0

6.3 Carbohydrates [i](#)

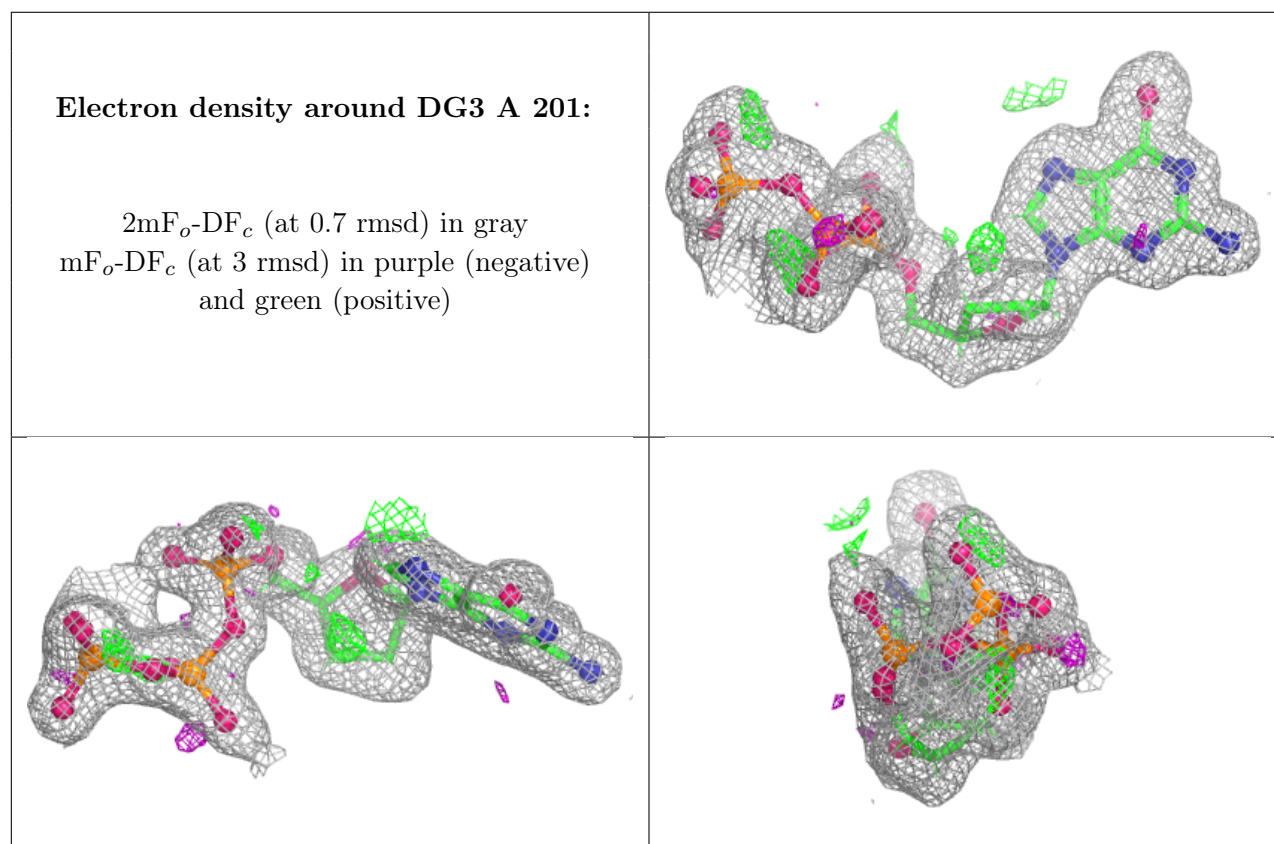
There are no oligosaccharides in this entry.

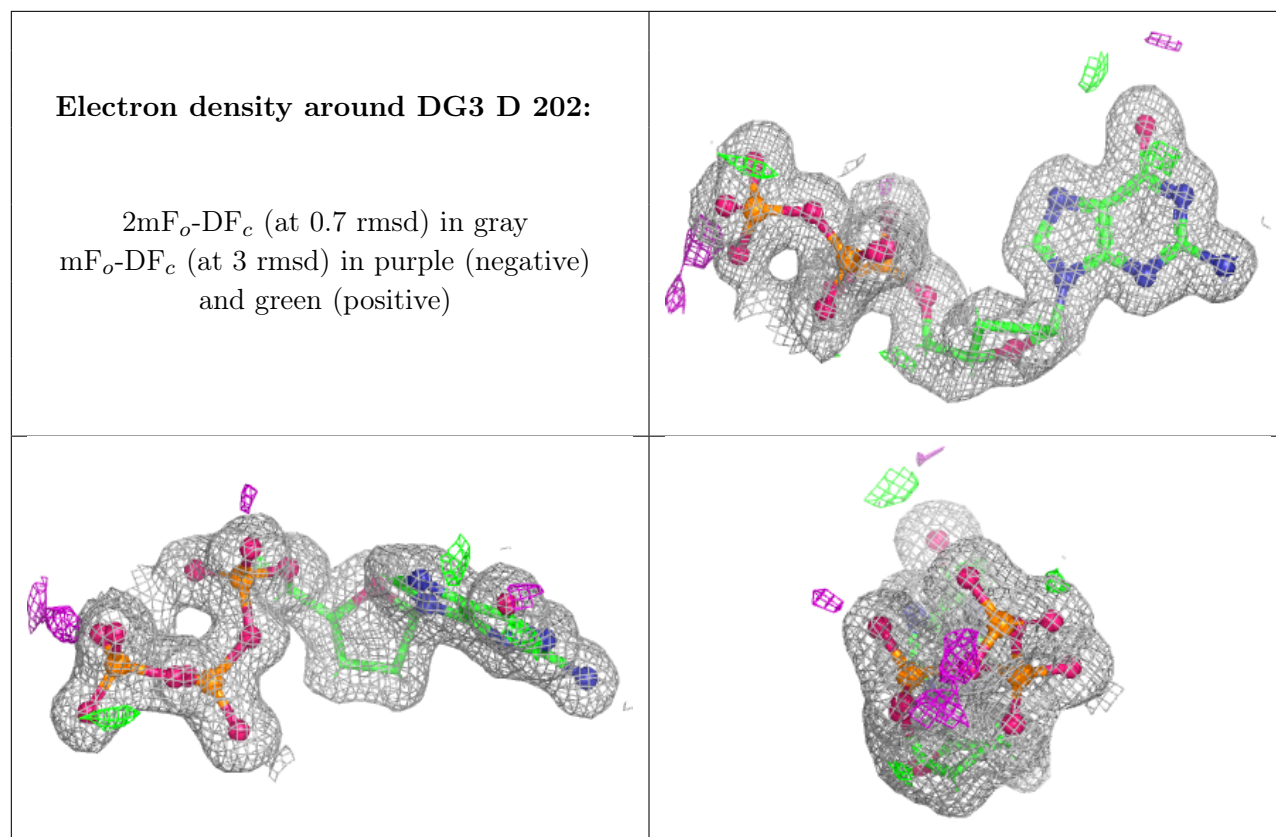
6.4 Ligands [i](#)

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($\text{\AA}^2$)	Q<0.9
6	SO4	A	2	5/5	0.92	0.08	35,47,51,52	0
6	SO4	D	1	5/5	0.93	0.11	28,36,45,47	0
4	DG3	A	201	30/30	0.95	0.08	16,21,28,33	0
5	MG	A	1	1/1	0.97	0.06	25,25,25,25	0
4	DG3	D	202	30/30	0.98	0.05	9,13,16,20	0
5	MG	D	2	1/1	0.99	0.03	13,13,13,13	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.