



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:56 PM UTC

PDB ID : 4EZ6 / pdb_00004ez6
Title : Bacillus DNA Polymerase I Large Fragment Complex 1
Authors : Wang, W.; Beese, L.S.
Deposited on : 2012-05-02
Resolution : 1.64 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

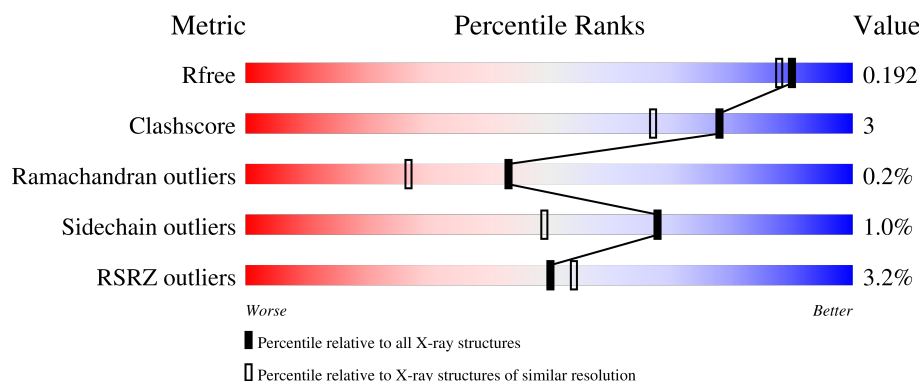
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.64 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	1141 (1.64-1.64)
Clashscore	190562	1171 (1.64-1.64)
Ramachandran outliers	187476	1151 (1.64-1.64)
Sidechain outliers	187428	1150 (1.64-1.64)
RSRZ outliers	180081	1141 (1.64-1.64)

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	4% 86% 6% 7%
1	D	592	% 92% 6% .
2	B	9	11% 67% 22% 11%
2	E	9	56% 22% 22%
3	C	13	8% 77% 8% 15%

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Mol	Chain	Length	Quality of chain
3	F	13	 A horizontal bar chart showing the quality of chain F. The bar is divided into three segments: a red segment on the left labeled '15%', a green segment in the middle labeled '92%', and a grey segment on the right labeled '8%'. The segments are separated by thin white lines.

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 21113 atoms, of which 9745 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.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	549	Total	C	H	N	O	S	0	5	0
			8949	2822	4524	770	818	15			
1	D	580	Total	C	H	N	O	S	0	5	0
			9424	2973	4750	811	873	17			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	598	ALA	ASP	engineered mutation	UNP Q5KWC1
A	710	TYR	PHE	engineered mutation	UNP Q5KWC1
A	823	HIS	ARG	SEE REMARK 999	UNP Q5KWC1
D	598	ALA	ASP	engineered mutation	UNP Q5KWC1
D	710	TYR	PHE	engineered mutation	UNP Q5KWC1
D	823	HIS	ARG	SEE REMARK 999	UNP Q5KWC1

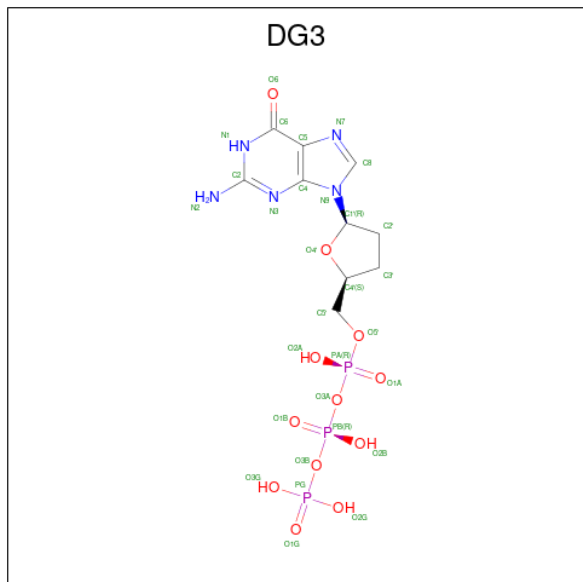
- Molecule 2 is a DNA chain called DNA (5'-D(*CP*CP*TP*GP*AP*CP*TP*CP*(DDG))-3').

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	9	Total	C	H	N	O	P	0	0	0
			277	86	100	31	52	8			
2	E	9	Total	C	H	N	O	P	0	0	0
			277	86	100	31	52	8			

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

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	C	11	Total	C	H	N	O	P	0	0	0
			353	108	123	45	66	11			
3	F	12	Total	C	H	N	O	P	0	0	0
			382	118	134	50	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
5	A	1	Total	C	H	O	0	0
			22	6	14	2		

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

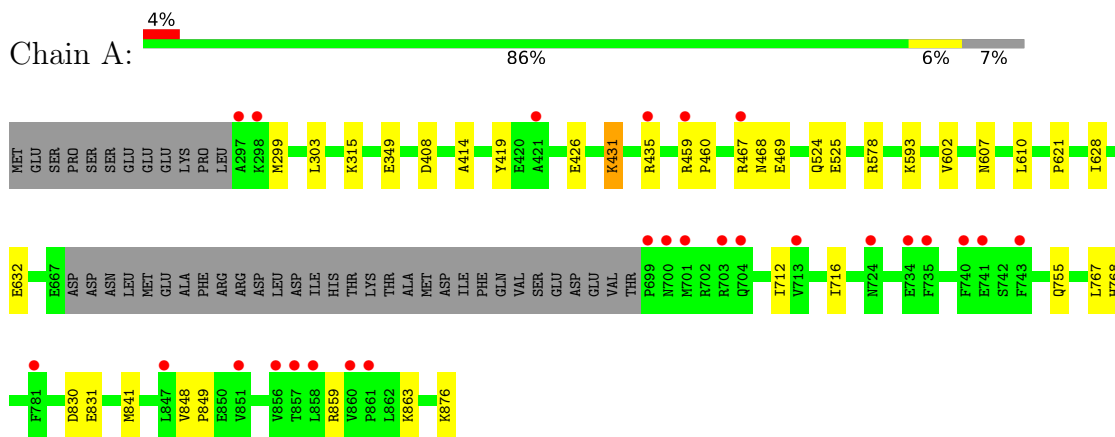
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	435	Total	O	0	0
			435	435		
7	D	719	Total	O	0	0
			719	719		
7	B	30	Total	O	0	0
			30	30		
7	C	68	Total	O	0	0
			68	68		
7	E	39	Total	O	0	0
			39	39		
7	F	73	Total	O	0	0
			73	73		

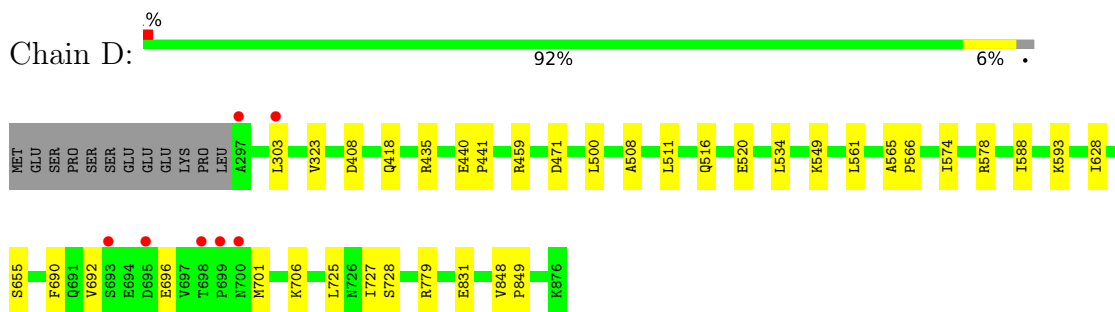
3 Residue-property plots [i](#)

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.

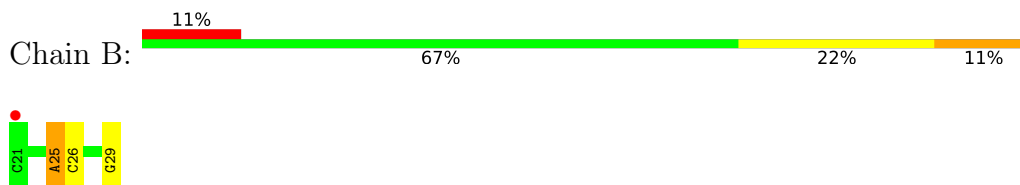
- Molecule 1: DNA polymerase



- Molecule 1: DNA polymerase

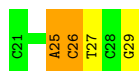


- Molecule 2: DNA (5'-D(*CP*CP*TP*GP*AP*CP*TP*CP*(DDG))-3')

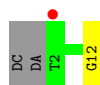
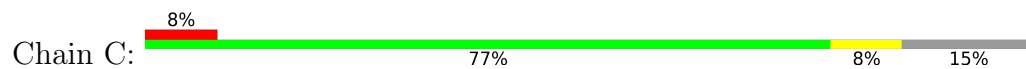


- Molecule 2: DNA (5'-D(*CP*CP*TP*GP*AP*CP*TP*CP*(DDG))-3')

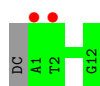
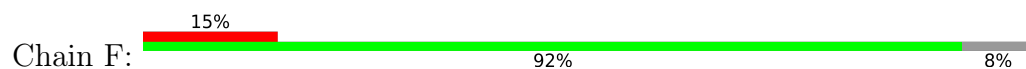




- Molecule 3: DNA (5'-D(*CP*AP*TP*GP*CP*GP*AP*GP*TP*CP*AP*GP*G)-3')



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	93.74Å 109.51Å 149.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.87 – 1.64 46.87 – 1.64	Depositor EDS
% Data completeness (in resolution range)	84.0 (46.87-1.64) 84.0 (46.87-1.64)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.49 (at 1.64Å)	Xtriage
Refinement program	PHENIX dev_1026	Depositor
R, R_{free}	0.171 , 0.192 0.171 , 0.192	Depositor DCC
R_{free} test set	7161 reflections (4.53%)	wwPDB-VP
Wilson B-factor (Å ²)	16.6	Xtriage
Anisotropy	0.229	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 49.5	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	21113	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.49% 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, MPD, 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	0.52	0/4520	0.73	2/6106 (0.0%)
1	D	0.70	0/4773	0.79	0/6452
2	B	0.51	0/173	1.34	3/264 (1.1%)
2	E	0.54	0/173	1.45	3/264 (1.1%)
3	C	0.44	0/258	1.04	0/397
3	F	0.57	0/279	1.21	0/430
All	All	0.61	0/10176	0.82	8/13913 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	25	DA	P-O5'-C5'	7.90	131.85	120.00
2	E	26	DC	N1-C1'-C2'	-7.27	102.59	113.50
2	E	25	DA	C5'-C4'-C3'	7.26	125.79	114.90
2	B	25	DA	P-O5'-C5'	6.33	129.49	120.00
2	B	25	DA	C5'-C4'-C3'	5.89	123.74	114.90
2	B	26	DC	N1-C1'-C2'	-5.55	105.17	113.50
1	A	859	ARG	N-CA-C	-5.46	106.62	113.28
1	A	299	MET	N-CA-C	5.10	117.10	108.02

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	4425	4524	4513	28	0
1	D	4674	4750	4738	24	0
2	B	177	100	103	1	0
2	E	177	100	103	3	0
3	C	230	123	124	1	0
3	F	248	134	136	0	0
4	A	30	0	12	0	0
4	D	30	0	12	2	0
5	A	8	14	14	1	0
6	D	5	0	0	0	0
7	A	435	0	0	15	0
7	B	30	0	0	0	0
7	C	68	0	0	1	0
7	D	719	0	0	7	1
7	E	39	0	0	0	0
7	F	73	0	0	0	0
All	All	11368	9745	9755	56	1

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 (56) 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:607[A]:ASN:OD1	7:A:1424:HOH:O	1.99	0.80
1:D:459:ARG:NH2	7:D:1486:HOH:O	2.16	0.78
5:A:902:MPD:HM2	7:A:1041:HOH:O	1.86	0.75
1:A:602[B]:VAL:HG21	1:A:621:PRO:HG3	1.68	0.74
1:A:876:LYS:OXT	7:A:1432:HOH:O	2.05	0.73
1:A:315:LYS:NZ	7:A:1433:HOH:O	2.22	0.71
1:A:459:ARG:NH2	7:A:1360:HOH:O	2.26	0.68
1:D:508:ALA:O	7:D:1377:HOH:O	2.13	0.66
1:D:831[B]:GLU:OE2	7:D:1185:HOH:O	2.14	0.63
1:A:830:ASP:OD2	7:A:1257:HOH:O	2.15	0.63
1:A:435:ARG:NE	7:A:1431:HOH:O	2.34	0.61
1:A:459:ARG:HB2	1:A:460:PRO:HD3	1.82	0.61
1:A:426:GLU:OE2	1:A:431:LYS:HD3	2.01	0.60
1:D:692:VAL:HG21	1:D:701:MET:HE1	1.85	0.59
1:A:632:GLU:HB2	7:A:1154:HOH:O	2.02	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:418:GLN:HA	7:D:1233:HOH:O	2.04	0.57
1:D:593:LYS:HG3	7:D:1442:HOH:O	2.03	0.57
3:C:12:DG:H1'	7:C:144:HOH:O	2.05	0.56
1:D:692:VAL:HB	1:D:696:GLU:HG3	1.88	0.56
1:A:435:ARG:NH2	7:A:1192:HOH:O	2.40	0.55
1:A:830:ASP:OD1	1:A:831:GLU:HG3	2.09	0.53
1:A:435:ARG:NH1	7:A:1316:HOH:O	2.33	0.53
1:D:323:VAL:O	1:D:435:ARG:NH2	2.43	0.52
1:D:500:LEU:HD11	1:D:588:ILE:HD13	1.91	0.51
1:D:779:ARG:NH1	7:D:1683:HOH:O	2.43	0.51
1:D:690:PHE:CD2	1:D:701:MET:HE3	2.46	0.50
1:A:593:LYS:NZ	7:A:1169:HOH:O	2.33	0.50
1:A:578:ARG:NH1	2:B:25:DA:H5''	2.28	0.48
1:A:467:ARG:NH2	7:A:1274:HOH:O	2.48	0.47
1:A:848:VAL:HB	1:A:849:PRO:HD3	1.95	0.47
1:A:767:LEU:O	1:A:768:HIS:HB2	2.14	0.47
1:A:435:ARG:NH2	7:A:1316:HOH:O	2.44	0.46
1:D:655:SER:HA	4:D:901:DG3:O2B	2.16	0.46
1:D:408:ASP:OD1	1:D:408:ASP:N	2.48	0.45
1:D:565:ALA:N	1:D:566:PRO:CD	2.80	0.45
1:A:408:ASP:OD1	1:A:408:ASP:N	2.47	0.45
1:D:534:LEU:HD11	1:D:574:ILE:HD13	1.98	0.44
1:D:706:LYS:HD3	4:D:901:DG3:H4'	2.00	0.44
1:A:468:ASN:O	1:A:469:GLU:HB2	2.18	0.44
1:D:848:VAL:HB	1:D:849:PRO:HD3	2.00	0.44
1:A:712:ILE:HA	1:A:716:ILE:HG22	2.00	0.43
2:E:26:DC:H2''	2:E:27:DT:O5'	2.17	0.43
1:A:435:ARG:CD	7:A:1431:HOH:O	2.67	0.43
1:D:435:ARG:NH2	7:D:1500:HOH:O	2.42	0.43
1:D:440:GLU:HB3	1:D:441:PRO:HD3	2.01	0.43
1:D:516:GLN:O	1:D:520:GLU:HG3	2.19	0.42
1:D:725:LEU:HB2	1:D:727:ILE:HG12	2.01	0.42
2:E:26:DC:H2'	2:E:27:DT:H71	2.00	0.42
1:A:459:ARG:CB	1:A:460:PRO:HD3	2.46	0.42
1:D:500:LEU:CD1	1:D:588:ILE:HD13	2.49	0.42
1:A:414:ALA:HB1	1:A:419:TYR:HB3	2.00	0.42
1:A:755:GLN:HG2	7:A:1201:HOH:O	2.19	0.41
1:D:578:ARG:NH1	2:E:25:DA:H5''	2.36	0.41
1:A:524:GLN:HG2	1:A:525:GLU:O	2.21	0.41
1:A:607[A]:ASN:CG	1:A:610:LEU:HB2	2.46	0.40
1:D:471:ASP:OD1	1:D:471:ASP:N	2.55	0.40

All (1) 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)
7:D:1240:HOH:O	7:D:1337:HOH:O[4_445]	2.08	0.12

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	550/592 (93%)	537 (98%)	12 (2%)	1 (0%)	43 27
1	D	583/592 (98%)	570 (98%)	12 (2%)	1 (0%)	43 27
All	All	1133/1184 (96%)	1107 (98%)	24 (2%)	2 (0%)	43 27

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	628	ILE
1	A	628	ILE

5.3.2 Protein sidechains [i](#)

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	471/507 (93%)	466 (99%)	5 (1%)	65 44
1	D	499/507 (98%)	494 (99%)	5 (1%)	68 49
All	All	970/1014 (96%)	960 (99%)	10 (1%)	68 49

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	303	LEU
1	A	349	GLU
1	A	431	LYS
1	A	841	MET
1	A	863	LYS
1	D	303	LEU
1	D	511	LEU
1	D	549	LYS
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	A	468	ASN
1	A	470	GLN
1	D	723	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	E	29	3,2	20,23,24	0.62	0	27,33,36	0.74	1 (3%)
2	DDG	B	29	3,2	20,23,24	0.66	0	27,33,36	0.60	1 (3%)

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	E	29	3,2	-	0/7/18/19	0/3/3/3
2	DDG	B	29	3,2	-	0/7/18/19	0/3/3/3

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	29	DDG	C4'-O4'-C1'	2.12	111.81	109.81
2	B	29	DDG	C2'-C3'-C4'	-2.01	99.01	102.75

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](#)

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	$\# Z > 2$
4	DG3	A	901	-	31,32,32	2.20	5 (16%)	44,50,50	1.85	11 (25%)
5	MPD	A	902	-	7,7,7	0.46	0	9,10,10	0.89	1 (11%)
4	DG3	D	901	-	31,32,32	2.16	9 (29%)	44,50,50	2.16	19 (43%)
6	SO4	D	902	-	4,4,4	0.27	0	6,6,6	0.46	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	DG3	A	901	-	-	6/22/31/31	0/3/3/3
5	MPD	A	902	-	-	4/5/5/5	-
4	DG3	D	901	-	-	3/22/31/31	0/3/3/3

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	901	DG3	O6-C6	8.62	1.39	1.23
4	D	901	DG3	O6-C6	6.31	1.35	1.23
4	D	901	DG3	PB-O3A	-5.97	1.53	1.59
4	A	901	DG3	C2-N2	4.90	1.45	1.34
4	D	901	DG3	C2-N2	3.52	1.42	1.34
4	A	901	DG3	PB-O3B	2.98	1.62	1.59
4	D	901	DG3	PA-O3A	-2.90	1.56	1.59
4	D	901	DG3	C6-N1	-2.46	1.34	1.38
4	D	901	DG3	PB-O3B	2.29	1.62	1.59
4	A	901	DG3	O4'-C1'	2.26	1.47	1.42
4	A	901	DG3	C2'-C3'	-2.16	1.48	1.54
4	D	901	DG3	C2'-C3'	-2.12	1.48	1.54
4	D	901	DG3	C2-N1	-2.09	1.32	1.37
4	D	901	DG3	C5-N7	-2.01	1.35	1.39

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	901	DG3	C5-C4-N3	-5.73	119.28	128.39
4	A	901	DG3	C5-C4-N3	-5.30	119.95	128.39
4	A	901	DG3	C2-N3-C4	4.52	120.08	112.30
4	D	901	DG3	C2-N3-C4	4.45	119.96	112.30
4	D	901	DG3	C2-N1-C6	-3.80	118.21	125.11
4	D	901	DG3	C4'-O4'-C1'	3.77	113.37	109.81
4	D	901	DG3	O5'-PA-O1A	3.55	123.02	108.94
4	D	901	DG3	N9-C4-N3	3.49	132.93	125.95
4	D	901	DG3	C5-C6-N1	3.29	121.62	113.25
4	A	901	DG3	C4-C5-N7	-3.09	105.77	110.67
4	A	901	DG3	C6-C5-N7	3.03	135.81	130.29
4	A	901	DG3	C2-N1-C6	-3.02	119.64	125.11
4	D	901	DG3	O3A-PA-O1A	-2.92	101.93	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	901	DG3	O4'-C1'-N9	2.85	112.91	107.86
4	A	901	DG3	O4'-C1'-N9	2.81	112.86	107.86
4	D	901	DG3	O4'-C4'-C5'	2.81	114.40	109.34
4	A	901	DG3	N9-C4-N3	2.76	131.48	125.95
4	A	901	DG3	C3'-C2'-C1'	2.73	106.02	102.87
4	D	901	DG3	O6-C6-C5	-2.65	119.54	126.53
4	D	901	DG3	C3'-C2'-C1'	2.48	105.73	102.87
4	D	901	DG3	O3G-PG-O3B	2.43	112.79	104.64
4	D	901	DG3	C2'-C1'-N9	2.41	116.97	112.40
4	D	901	DG3	O2A-PA-O3A	-2.39	100.81	107.27
4	A	901	DG3	C5-C6-N1	2.34	119.22	113.25
4	A	901	DG3	C8-N7-C5	2.34	108.43	104.26
4	D	901	DG3	C6-C5-N7	2.31	134.50	130.29
4	D	901	DG3	O4'-C1'-C2'	-2.29	103.70	106.41
5	A	902	MPD	O2-C2-C3	2.20	117.48	109.27
4	A	901	DG3	C4'-O4'-C1'	2.06	111.75	109.81
4	D	901	DG3	C4-C5-N7	-2.04	107.43	110.67
4	D	901	DG3	PA-O5'-C5'	2.01	132.84	121.35

There are no chirality outliers.

All (13) torsion outliers are listed below:

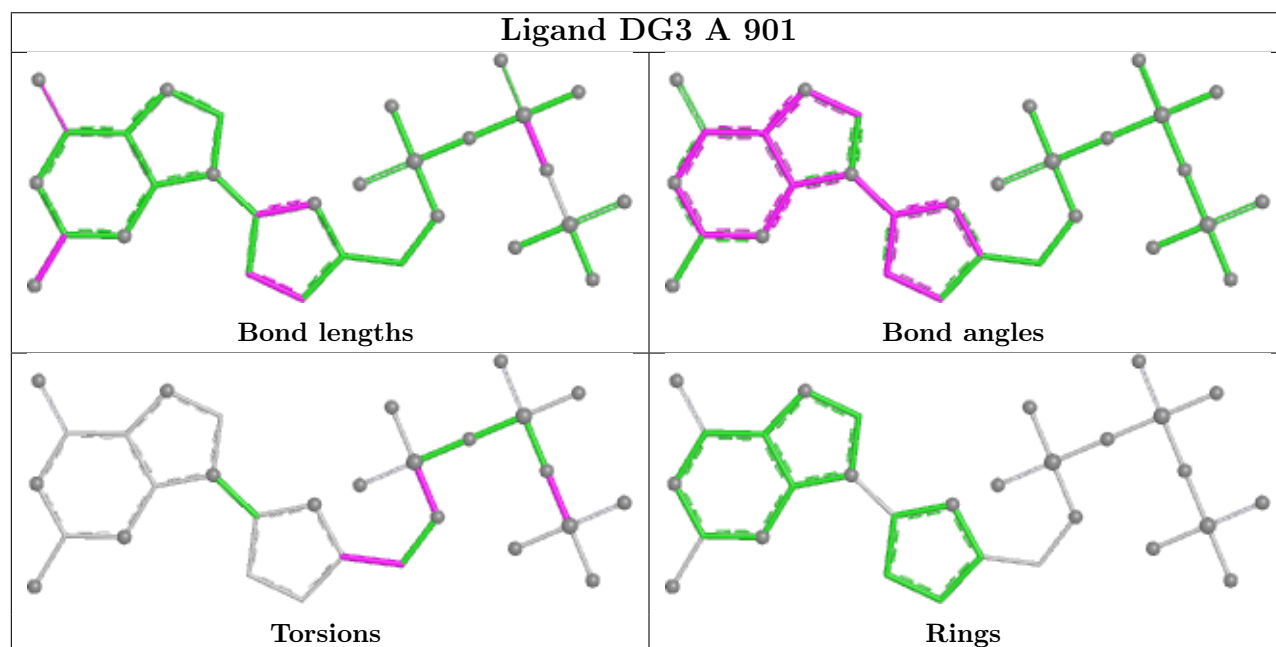
Mol	Chain	Res	Type	Atoms
4	A	901	DG3	C5'-O5'-PA-O3A
4	A	901	DG3	O4'-C4'-C5'-O5'
4	A	901	DG3	C3'-C4'-C5'-O5'
4	D	901	DG3	C3'-C4'-C5'-O5'
4	A	901	DG3	PB-O3B-PG-O3G
5	A	902	MPD	C2-C3-C4-O4
5	A	902	MPD	CM-C2-C3-C4
4	A	901	DG3	C5'-O5'-PA-O1A
4	D	901	DG3	O4'-C1'-N9-C4
4	D	901	DG3	O4'-C4'-C5'-O5'
4	A	901	DG3	PB-O3B-PG-O2G
5	A	902	MPD	O2-C2-C3-C4
5	A	902	MPD	C1-C2-C3-C4

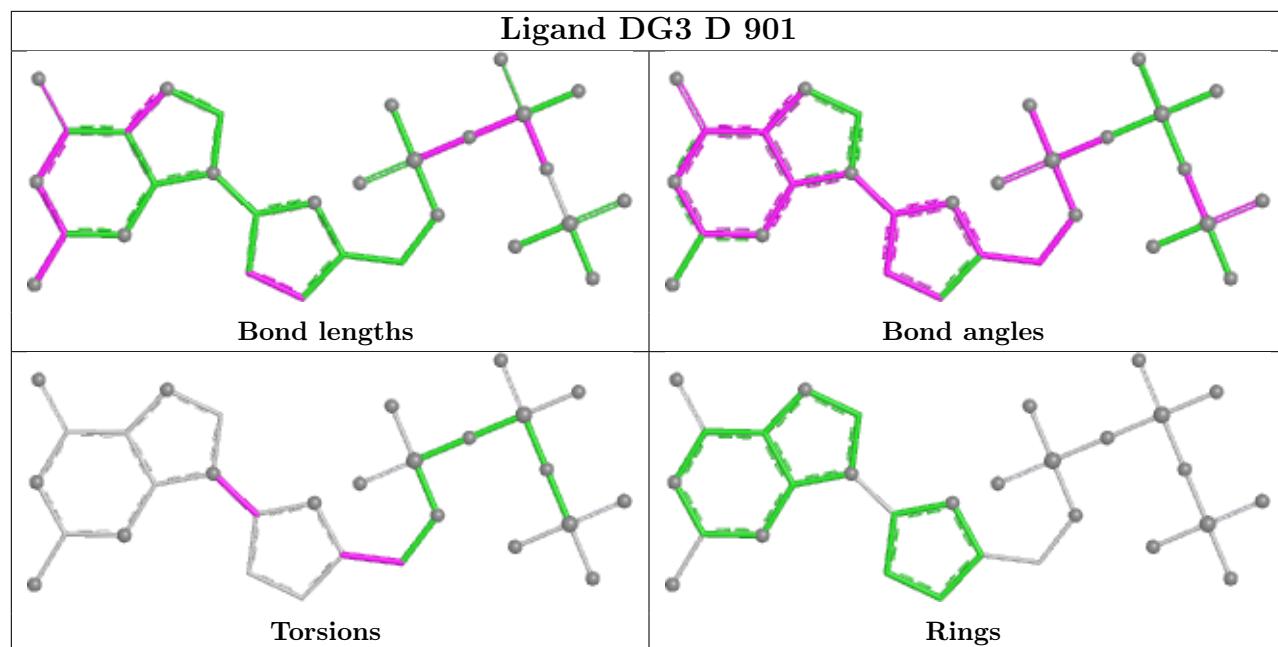
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	902	MPD	1	0
4	D	901	DG3	2	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	549/592 (92%)	0.42	26 (4%) 36 41	11, 36, 61, 82	5 (0%)
1	D	580/592 (97%)	-0.27	7 (1%) 76 81	7, 20, 41, 61	5 (0%)
2	B	8/9 (88%)	0.11	1 (12%) 8 9	21, 26, 46, 60	0
2	E	8/9 (88%)	-0.22	0 100 100	13, 21, 43, 51	0
3	C	11/13 (84%)	-0.15	1 (9%) 15 17	18, 26, 41, 64	0
3	F	12/13 (92%)	-0.19	2 (16%) 4 5	11, 21, 54, 62	0
All	All	1168/1228 (95%)	0.06	37 (3%) 50 54	7, 28, 56, 82	10 (0%)

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	297	ALA	4.5
1	A	699	PRO	4.2
1	A	857	THR	3.7
1	A	735	PHE	3.5
1	D	698	THR	3.5
1	A	860	VAL	3.5
1	D	699	PRO	3.2
1	D	693	SER	3.1
1	A	847	LEU	3.0
1	A	858	LEU	3.0
1	A	734	GLU	3.0
3	F	1	DA	2.9
1	A	851	VAL	2.9
1	D	297	ALA	2.9
1	A	861	PRO	2.7
1	A	298	LYS	2.5
1	A	701	MET	2.5
1	A	740	PHE	2.4
1	D	303	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	741	GLU	2.4
1	A	467	ARG	2.4
1	A	781	PHE	2.4
1	D	695	ASP	2.4
1	A	743	PHE	2.3
1	A	703	ARG	2.3
1	D	700	ASN	2.3
3	C	2	DT	2.2
1	A	724	ASN	2.2
3	F	2	DT	2.2
1	A	704	GLN	2.2
1	A	713	VAL	2.2
1	A	421	ALA	2.2
1	A	700	ASN	2.2
1	A	856	VAL	2.1
1	A	459	ARG	2.1
2	B	21	DC	2.1
1	A	435	ARG	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	DDG	B	29	21/22	0.97	0.06	21,25,30,34	0
2	DDG	E	29	21/22	0.99	0.04	10,13,17,18	0

6.3 Carbohydrates [i](#)

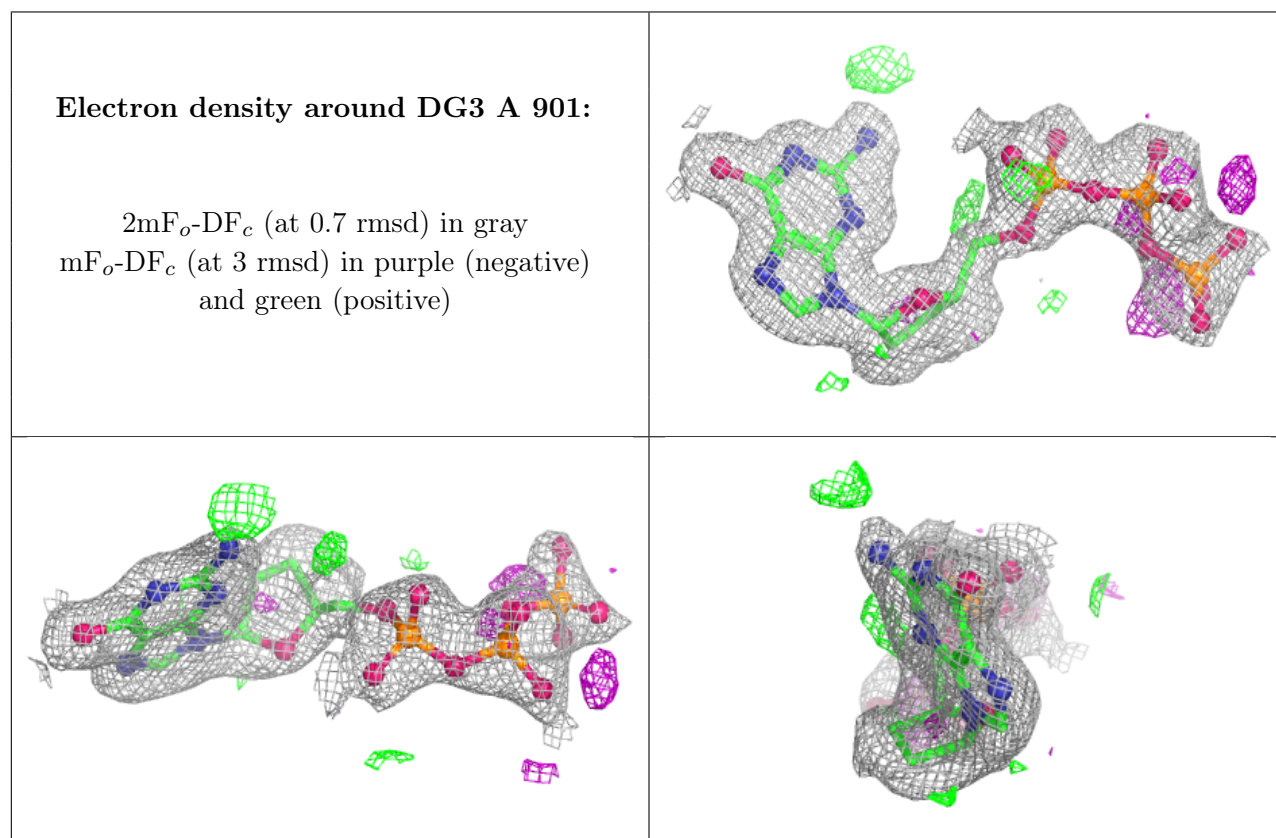
There are no oligosaccharides in this entry.

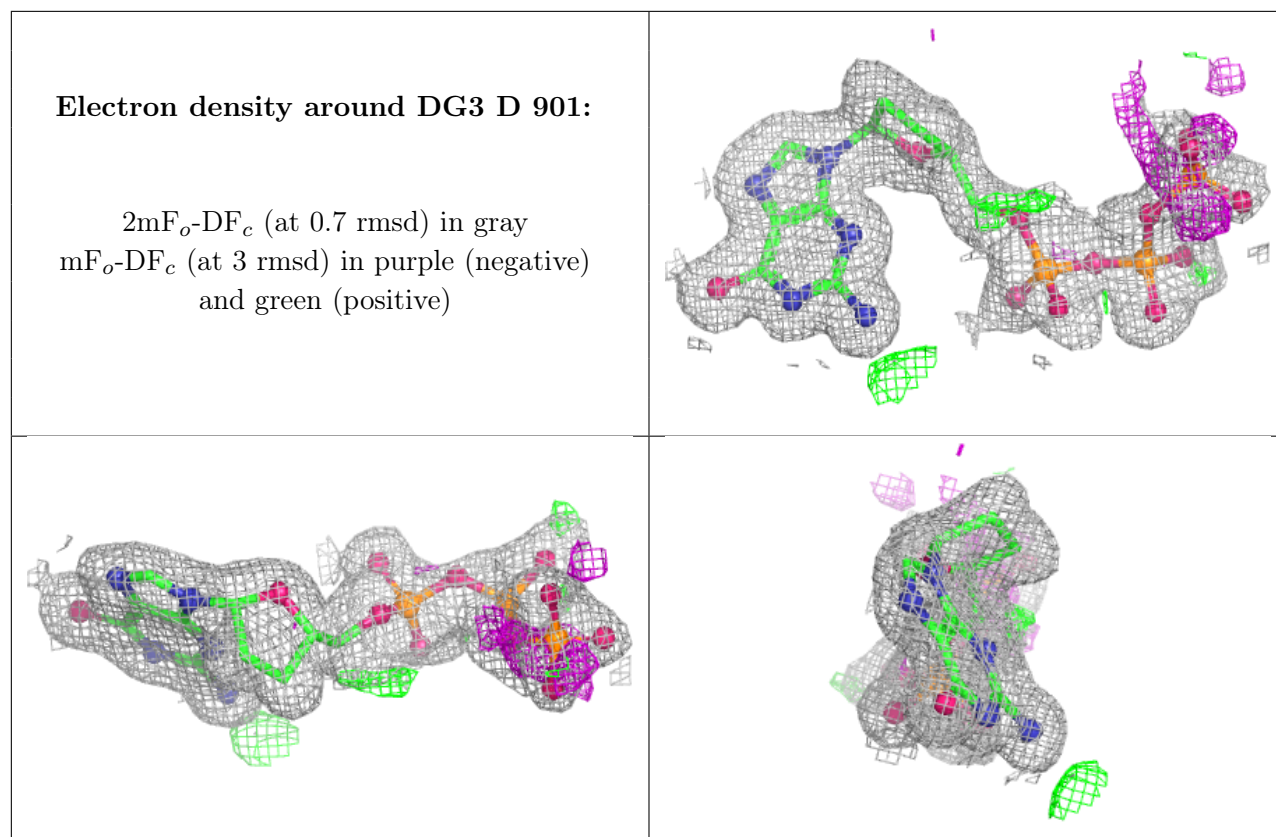
6.4 Ligands [i](#)

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
5	MPD	A	902	8/8	0.89	0.10	17,28,34,41	0
4	DG3	A	901	30/30	0.93	0.10	18,27,66,85	0
6	SO4	D	902	5/5	0.94	0.09	28,33,42,49	0
4	DG3	D	901	30/30	0.96	0.09	11,19,42,49	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.