



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:45 PM UTC

PDB ID : 4KHN / pdb_00004khn
Title : Crystal structure of the ternary complex of the D714A mutant of RB69 DNA polymerase
Authors : Guja, K.E.; Jacewicz, A.; Trzemecka, A.; Plochocka, D.; Yakubovskaya, E.; Bebenek, A.; Garcia-Diaz, M.
Deposited on : 2013-04-30
Resolution : 2.55 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

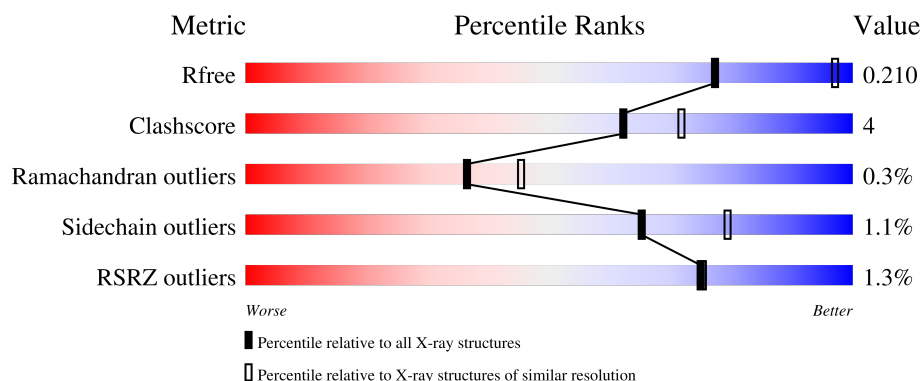
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	903	1% 91% 9%
1	B	903	2% 82% 13% ••
2	C	18	94% 6%
2	E	18	94% 6%

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Mol	Chain	Length	Quality of chain
3	D	13	 100%
3	F	13	 92%  8%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
7	GOL	A	1008	-	-	X	-

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 15657 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	901	Total	C	N	O	S	0	2	0
			7198	4640	1199	1326	33			
1	B	870	Total	C	N	O	S	0	1	0
			6730	4352	1113	1235	30			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	222	ALA	ASP	engineered mutation	UNP Q38087
A	327	ALA	ASP	engineered mutation	UNP Q38087
A	567	ALA	TYR	engineered mutation	UNP Q38087
A	714	ALA	ASP	engineered mutation	UNP Q38087
B	222	ALA	ASP	engineered mutation	UNP Q38087
B	327	ALA	ASP	engineered mutation	UNP Q38087
B	567	ALA	TYR	engineered mutation	UNP Q38087
B	714	ALA	ASP	engineered mutation	UNP Q38087

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	18	Total	C	N	O	P	0	0	0
			365	174	69	105	17			
2	E	18	Total	C	N	O	P	0	0	0
			365	174	69	105	17			

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

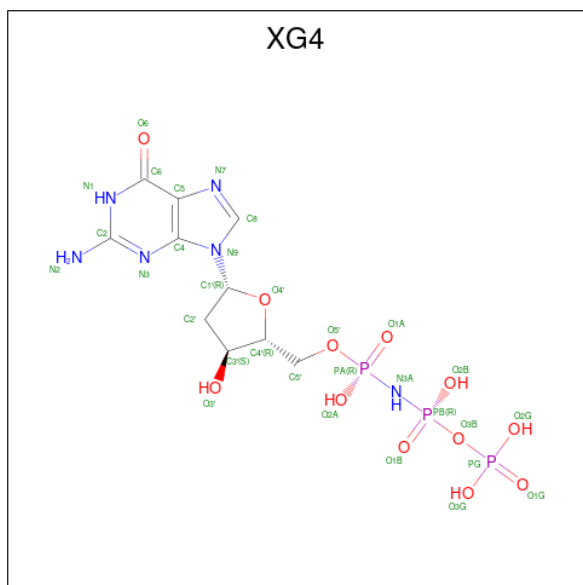
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	13	Total	C	N	O	P	0	0	0
			263	126	48	77	12			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	13	Total	C	N	O	P	0	0	0
			263	126	48	77	12			

- Molecule 4 is 2'-deoxy-5'-O-[(R)-hydroxy{[(R)-hydroxy(phosphonooxy)phosphoryl]amino}phosphoryl]guanosine (CCD ID: XG4) (formula: C₁₀H₁₇N₆O₁₂P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			31	10	6	12	3		
4	C	1	Total	C	N	O	P	0	0
			31	10	6	12	3		
4	B	1	Total	C	N	O	P	0	0
			31	10	6	12	3		
4	B	1	Total	C	N	O	P	0	0
			31	10	6	12	3		

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Ca 1 1	0	0
5	B	1	Total Ca 1 1	0	0

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



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

- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 8 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	1	Total	Na	0	0
			1	1		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	228	Total	O	0	0
			228	228		
9	C	23	Total	O	0	0
			23	23		
9	D	17	Total	O	0	0
			17	17		
9	E	6	Total	O	0	0
			6	6		
9	F	3	Total	O	0	0
			3	3		
9	B	27	Total	O	0	0
			27	27		

Chain C:  94% 6%

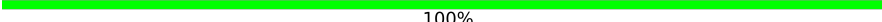


- Molecule 2: DNA (5'-D(*TP*CP*AP*CP*GP*TP*AP*AP*GP*CP*AP*GP*TP*CP*CP*GP*CP*G)-3')

Chain E:  94% 6%



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

Chain D:  100%

There are no outlier residues recorded for this chain.

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

Chain F:  92% 8%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.26Å 119.35Å 148.03Å 90.00° 91.64° 90.00°	Depositor
Resolution (Å)	35.33 – 2.55 35.33 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.8 (35.33-2.55) 95.4 (35.33-2.55)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.27 (at 2.54Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1363)	Depositor
R, R_{free}	0.168 , 0.211 0.171 , 0.210	Depositor DCC
R_{free} test set	4149 reflections (3.70%)	wwPDB-VP
Wilson B-factor (Å ²)	51.6	Xtriage
Anisotropy	0.236	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 54.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.032 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	15657	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.58% 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: CA, NA, GOL, XG4, 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	0.59	0/7382	0.88	3/9996 (0.0%)
1	B	0.55	1/6904 (0.0%)	0.90	10/9388 (0.1%)
2	C	0.13	0/409	0.55	0/629
2	E	0.13	0/409	0.55	0/629
3	D	0.12	0/294	0.61	0/452
3	F	0.09	0/294	0.59	0/452
All	All	0.55	1/15692 (0.0%)	0.86	13/21546 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	821	ALA	C-N	5.02	1.40	1.33

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	355	ILE	N-CA-C	6.93	123.75	109.34
1	B	798	GLY	CA-C-N	6.38	126.33	119.76
1	B	798	GLY	C-N-CA	6.38	126.33	119.76
1	B	821	ALA	CA-C-N	-6.36	113.21	119.76
1	B	821	ALA	C-N-CA	-6.36	113.21	119.76
1	A	660	GLU	CA-C-N	-6.01	112.38	119.05
1	A	660	GLU	C-N-CA	-6.01	112.38	119.05
1	A	900	MET	CB-CG-SD	-5.95	94.86	112.70
1	B	85	MET	N-CA-C	5.85	117.74	108.67
1	B	896	SER	N-CA-C	-5.53	100.66	109.07
1	B	101	ILE	N-CA-C	5.29	114.88	107.37
1	B	243	SER	CA-C-N	5.03	124.69	119.56
1	B	243	SER	C-N-CA	5.03	124.69	119.56

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	7198	0	6994	42	0
1	B	6730	0	6330	74	0
2	C	365	0	203	1	0
2	E	365	0	203	1	0
3	D	263	0	148	0	0
3	F	263	0	148	1	0
4	A	31	0	13	0	0
4	B	62	0	26	2	0
4	C	31	0	13	2	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	25	0	0	0	0
6	B	5	0	0	0	0
7	A	12	0	16	4	0
8	A	1	0	0	0	0
9	A	228	0	0	1	0
9	B	27	0	0	0	0
9	C	23	0	0	0	0
9	D	17	0	0	0	0
9	E	6	0	0	0	0
9	F	3	0	0	0	0
All	All	15657	0	14094	121	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 4.

All (121) 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:395:PHE:HB2	1:B:591:GLN:HG3	1.61	0.82
1:B:560:LYS:NZ	4:B:1002:XG4:O3G	2.22	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:153:ASN:ND2	1:B:192:ASP:O	2.22	0.71
1:A:833:LEU:HD13	1:A:866:MET:HG2	1.73	0.70
1:A:514:LEU:HD21	1:A:529:LYS:HG2	1.74	0.69
1:B:546:GLN:HE21	1:B:550:VAL:HG23	1.59	0.68
1:A:664:ASP:OD2	1:A:668:ARG:NH1	2.26	0.68
1:A:412:LEU:HG	1:A:683:MET:HE3	1.77	0.67
7:A:1008:GOL:H31	4:C:101:XG4:HN3A	1.59	0.67
1:A:752:MET:HE3	1:A:889:LEU:HD12	1.79	0.65
1:A:486:LYS:HE2	1:A:490:LEU:HD11	1.79	0.64
1:B:170:LEU:HA	1:B:177:GLU:HG3	1.80	0.63
1:A:820:ASP:OD1	1:A:820:ASP:N	2.33	0.62
1:A:668:ARG:HD2	9:A:1290:HOH:O	1.98	0.61
1:B:417:PRO:HG2	1:B:563:ILE:HD12	1.83	0.61
7:A:1008:GOL:O3	7:A:1008:GOL:O1	2.18	0.60
1:A:395:PHE:HB2	1:A:591:GLN:HB2	1.86	0.58
1:B:475:ILE:HD12	1:B:566:LEU:HD23	1.86	0.58
1:A:303:LEU:HD21	1:A:319:ARG:HG3	1.86	0.57
1:B:412:LEU:HG	1:B:683:MET:HE3	1.86	0.57
1:B:35:PRO:HG3	1:B:65:MET:HG2	1.86	0.55
1:B:36:SER:OG	1:B:59:ARG:NH1	2.39	0.55
1:B:471:VAL:HG13	1:B:566:LEU:HD21	1.88	0.54
1:B:41:CYS:HB3	1:B:58:THR:HG22	1.90	0.54
1:B:209:THR:HG21	1:B:244:PRO:HG3	1.89	0.53
1:B:606:ASN:OD1	1:B:613:GLY:N	2.42	0.53
1:A:186:ILE:HD13	1:A:325:ILE:HD13	1.91	0.53
1:B:367:ALA:O	1:B:371:ASN:ND2	2.32	0.53
1:B:712:VAL:HG11	1:B:715:MET:HE2	1.90	0.53
1:B:109:ARG:HD2	1:B:140:ASP:OD1	2.09	0.52
1:A:900:MET:HE3	1:A:901:PHE:H	1.74	0.52
1:B:15:ILE:HD13	1:B:92:TYR:CE2	2.45	0.52
1:A:303:LEU:HD23	1:A:323:TYR:HB2	1.90	0.52
1:A:347[A]:MET:HE3	1:A:562:LEU:HD13	1.91	0.51
1:A:36:SER:OG	1:A:59:ARG:NH1	2.42	0.51
1:B:291:ASP:OD1	1:B:302:LYS:HA	2.10	0.51
1:B:226:VAL:O	1:B:230:ILE:HG12	2.10	0.51
1:A:436:VAL:HG12	1:A:437:ALA:O	2.10	0.51
1:B:50:PHE:HD1	1:B:56:PRO:HA	1.77	0.50
1:B:105:HIS:CE1	1:B:106:THR:HG23	2.47	0.50
1:B:482:ARG:HB2	1:B:559:ARG:HB3	1.92	0.50
1:A:465:LYS:NZ	1:A:675:ASN:OD1	2.44	0.49
1:B:145:ARG:HH11	1:B:185:LYS:HA	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:436:VAL:HG12	1:B:437:ALA:O	2.12	0.49
1:B:116:GLU:HB2	1:B:135:ALA:HB3	1.96	0.48
1:A:787:ASN:HB3	1:A:790:LYS:HB3	1.96	0.47
1:A:133:ILE:HG13	1:A:229:ARG:HG2	1.97	0.47
1:B:50:PHE:CD1	1:B:56:PRO:HA	2.50	0.47
1:B:482:ARG:NE	1:B:560:LYS:HB2	2.30	0.47
1:A:441:ASP:HB3	1:A:447:ALA:HB2	1.95	0.47
1:B:555:ALA:HB1	1:B:559:ARG:NH1	2.29	0.47
1:B:659:MET:O	1:B:663:ILE:HG13	2.15	0.47
1:A:540:GLU:HG3	1:A:544:ARG:HE	1.79	0.47
1:B:553:MET:O	1:B:556:GLN:HG3	2.13	0.47
1:B:731:GLU:OE1	1:B:731:GLU:N	2.44	0.47
1:B:109:ARG:HG2	1:B:210:PRO:HA	1.97	0.47
1:B:353:ILE:CD1	1:B:357:SER:HB2	2.45	0.47
4:B:1002:XG4:O2G	4:B:1002:XG4:O2B	2.33	0.46
1:A:706:LYS:HE3	2:C:7:DA:N3	2.31	0.46
1:B:752:MET:HE3	1:B:889:LEU:HD12	1.98	0.46
1:A:471:VAL:HG13	1:A:566:LEU:HD21	1.97	0.46
1:A:771:PHE:CZ	1:A:872:LEU:HB2	2.51	0.46
1:B:772:ARG:HG2	1:B:868:TYR:CG	2.51	0.46
1:B:116:GLU:HG2	1:B:324:ASN:ND2	2.31	0.45
1:B:546:GLN:NE2	1:B:550:VAL:HG23	2.27	0.45
1:B:752:MET:HG2	1:B:760:LEU:HD22	1.97	0.45
1:B:458:PRO:HG2	1:B:588:THR:HG22	1.98	0.45
1:B:438:PRO:HG2	1:B:441:ASP:OD2	2.16	0.45
1:B:777:ILE:HD13	1:B:848:TRP:HZ2	1.81	0.45
1:B:441:ASP:HB3	1:B:447:ALA:HB2	1.98	0.45
2:E:8:DA:H5"	1:B:705:LYS:HD3	1.98	0.44
1:A:380:ILE:HD12	1:A:576:ARG:CZ	2.48	0.44
1:A:353:ILE:HD12	1:A:357:SER:HB2	1.99	0.44
1:A:854:ILE:HD13	1:A:862:VAL:HG21	2.00	0.44
1:B:39:ALA:HB2	1:B:71:TRP:HH2	1.83	0.44
1:B:499:ILE:HD12	1:B:538:LEU:HD22	1.99	0.44
1:B:553:MET:O	1:B:557:ILE:HG13	2.18	0.44
1:B:597:ILE:HD12	1:B:597:ILE:HA	1.76	0.44
1:B:353:ILE:HD13	1:B:357:SER:HB2	2.00	0.43
1:B:126:PRO:HB2	1:B:224:PRO:HB2	2.00	0.43
1:B:273:TYR:OH	1:B:335:ASP:HA	2.19	0.43
1:B:499:ILE:HD13	1:B:499:ILE:HA	1.85	0.43
1:A:862:VAL:O	1:A:866:MET:HG3	2.19	0.43
1:A:179:PRO:HG2	1:A:329:TYR:CE2	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:49:TYR:CZ	1:B:59:ARG:HD3	2.54	0.43
1:B:352:LYS:HE3	1:B:352:LYS:HB2	1.63	0.43
1:B:757:GLU:HB2	1:B:889:LEU:HD22	2.01	0.43
1:B:878:LYS:HB3	1:B:879:PRO:HD3	2.00	0.43
3:F:113:DT:H5'	1:B:734:LYS:HG2	2.00	0.43
1:B:3:GLU:HG3	1:B:21:ASP:HA	2.00	0.43
1:B:594:LEU:HD12	1:B:594:LEU:HA	1.86	0.43
1:B:881:GLU:O	1:B:885:SER:HB3	2.18	0.43
7:A:1008:GOL:H32	4:C:101:XG4:N7	2.34	0.42
1:B:126:PRO:O	1:B:228:ASN:ND2	2.52	0.42
1:B:191:PHE:CZ	1:B:200:GLU:HG2	2.54	0.42
1:B:270:VAL:O	1:B:271:LEU:HD23	2.19	0.42
1:B:15:ILE:O	1:B:15:ILE:HD12	2.19	0.42
7:A:1008:GOL:HO1	7:A:1008:GOL:HO3	1.61	0.42
1:A:731:GLU:OE1	1:A:731:GLU:N	2.48	0.42
1:A:737:THR:HA	1:A:738:PRO:HD3	1.89	0.42
1:A:214:THR:OG1	1:A:215:GLY:N	2.52	0.42
1:B:162:TRP:CD1	1:B:321:ILE:HB	2.55	0.42
1:A:381:PRO:O	1:A:576:ARG:HD3	2.20	0.42
1:B:229:ARG:O	1:B:233:ILE:HG12	2.19	0.42
1:A:162:TRP:CH2	1:A:164:ILE:HG13	2.55	0.41
1:A:878:LYS:HB3	1:A:879:PRO:HD3	2.02	0.41
1:B:89:LYS:O	1:B:93:LEU:HG	2.20	0.41
1:B:285:GLN:HA	1:B:286:PRO:HD3	1.89	0.41
1:A:125:GLU:HA	1:A:126:PRO:HD2	1.90	0.41
1:A:44:SER:C	1:A:46:ALA:H	2.29	0.41
1:A:204:PHE:CE1	1:A:208:LYS:HD2	2.56	0.41
1:A:373:LEU:HD12	1:A:380:ILE:HG22	2.03	0.41
1:B:364:THR:O	1:B:368:ILE:HG13	2.20	0.41
1:B:226:VAL:HG12	1:B:242:LEU:HD11	2.02	0.41
1:A:302:LYS:HE2	1:A:323:TYR:CZ	2.55	0.41
1:B:555:ALA:HB1	1:B:559:ARG:HH12	1.85	0.41
1:B:580:LEU:HD12	1:B:580:LEU:HA	1.81	0.41
1:A:243:SER:HA	1:A:244:PRO:HD2	1.83	0.40
1:B:11:ILE:HB	1:B:16:PHE:CE1	2.56	0.40
1:A:805:ILE:HD13	1:A:808:ILE:HD12	2.03	0.40
1:B:348:GLY:HA3	1:B:355:ILE:HD13	2.04	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	901/903 (100%)	870 (97%)	28 (3%)	3 (0%)	36	45
1	B	867/903 (96%)	832 (96%)	32 (4%)	3 (0%)	36	45
All	All	1768/1806 (98%)	1702 (96%)	60 (3%)	6 (0%)	36	45

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	355	ILE
1	B	414	SER
1	B	819	ILE
1	A	414	SER
1	A	622	THR
1	A	424	ASN

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	749/798 (94%)	744 (99%)	5 (1%)	76	85
1	B	663/798 (83%)	652 (98%)	11 (2%)	53	72
All	All	1412/1596 (88%)	1396 (99%)	16 (1%)	65	79

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	ARG
1	A	557	ILE
1	A	608	VAL
1	A	624	SER
1	A	819	ILE
1	B	15	ILE
1	B	28	THR
1	B	177	GLU
1	B	180	SER
1	B	352	LYS
1	B	353	ILE
1	B	435	LYS
1	B	499	ILE
1	B	536	LYS
1	B	580	LEU
1	B	618	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	153	ASN
1	A	158	ASN
1	A	485	HIS
1	A	546	GLN
1	A	558	ASN
1	A	733	GLN
1	A	787	ASN
1	A	818	ASN
1	B	10	GLN
1	B	105	HIS
1	B	217	ASN
1	B	546	GLN
1	B	733	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 3 are monoatomic - leaving 12 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	SO4	B	1003	-	4,4,4	0.23	0	6,6,6	0.12	0
4	XG4	B	1001	-	32,33,33	1.71	4 (12%)	46,52,52	1.85	8 (17%)
4	XG4	C	101	5	32,33,33	1.53	5 (15%)	46,52,52	2.07	12 (26%)
4	XG4	A	1001	-	32,33,33	1.75	5 (15%)	46,52,52	1.80	6 (13%)
6	SO4	A	1007	-	4,4,4	0.24	0	6,6,6	0.07	0
6	SO4	A	1003	-	4,4,4	0.27	0	6,6,6	0.11	0
6	SO4	A	1004	-	4,4,4	0.23	0	6,6,6	0.18	0
6	SO4	A	1005	-	4,4,4	0.25	0	6,6,6	0.12	0
4	XG4	B	1002	5	32,33,33	1.65	6 (18%)	46,52,52	1.86	8 (17%)
7	GOL	A	1008	-	5,5,5	0.23	0	5,5,5	1.03	0
7	GOL	A	1009	-	5,5,5	0.41	0	5,5,5	0.30	0
6	SO4	A	1006	-	4,4,4	0.22	0	6,6,6	0.15	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	XG4	B	1001	-	-	10/19/34/34	0/3/3/3
4	XG4	C	101	5	-	9/19/34/34	0/3/3/3
4	XG4	A	1001	-	-	5/19/34/34	0/3/3/3
4	XG4	B	1002	5	-	8/19/34/34	0/3/3/3
7	GOL	A	1008	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	GOL	A	1009	-	-	2/4/4/4	-

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1001	XG4	PA-N3A	4.71	1.75	1.63
4	A	1001	XG4	PA-N3A	4.68	1.75	1.63
4	A	1001	XG4	PB-N3A	4.65	1.75	1.63
4	B	1001	XG4	PB-N3A	4.63	1.75	1.63
4	B	1002	XG4	PB-N3A	4.50	1.75	1.63
4	C	101	XG4	PA-N3A	4.04	1.73	1.63
4	B	1002	XG4	PA-N3A	4.00	1.73	1.63
4	C	101	XG4	PB-N3A	3.60	1.72	1.63
4	B	1001	XG4	C5-C4	3.16	1.47	1.38
4	B	1002	XG4	C5-C4	3.16	1.47	1.38
4	C	101	XG4	C5-C4	3.11	1.47	1.38
4	A	1001	XG4	C5-C4	3.09	1.47	1.38
4	A	1001	XG4	C6-N1	-3.07	1.33	1.38
4	C	101	XG4	C6-N1	-2.52	1.34	1.38
4	B	1002	XG4	C6-N1	-2.45	1.34	1.38
4	B	1001	XG4	C6-N1	-2.31	1.34	1.38
4	B	1002	XG4	C5-N7	-2.27	1.34	1.39
4	C	101	XG4	C5-N7	-2.24	1.34	1.39
4	B	1002	XG4	PB-O3B	2.22	1.61	1.59
4	A	1001	XG4	C5-N7	-2.08	1.34	1.39

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1001	XG4	C5-C4-N3	-6.48	118.08	128.39
4	C	101	XG4	C5-C4-N3	-6.36	118.27	128.39
4	A	1001	XG4	C5-C4-N3	-6.21	118.51	128.39
4	B	1002	XG4	C5-C4-N3	-6.02	118.81	128.39
4	B	1001	XG4	C2-N3-C4	5.36	121.53	112.30
4	C	101	XG4	C2-N3-C4	5.24	121.32	112.30
4	A	1001	XG4	N9-C4-N3	5.06	136.06	125.95
4	B	1001	XG4	N9-C4-N3	5.02	135.99	125.95
4	A	1001	XG4	C2-N3-C4	4.92	120.78	112.30
4	B	1002	XG4	C2-N3-C4	4.83	120.62	112.30
4	C	101	XG4	O1A-PA-N3A	-4.58	105.03	111.77
4	B	1002	XG4	N9-C4-N3	4.51	134.97	125.95
4	C	101	XG4	N9-C4-N3	4.46	134.86	125.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1002	XG4	C2'-C1'-N9	-3.42	105.25	113.81
4	C	101	XG4	C2'-C1'-N9	-3.33	105.48	113.81
4	C	101	XG4	C6-C5-N7	3.21	136.14	130.29
4	B	1002	XG4	C6-C5-N7	2.93	135.61	130.29
4	A	1001	XG4	C6-C5-N7	2.91	135.58	130.29
4	B	1001	XG4	C6-C5-N7	2.86	135.49	130.29
4	B	1002	XG4	O6-C6-C5	-2.60	119.68	126.53
4	C	101	XG4	C4-C5-N7	-2.57	106.60	110.67
4	B	1001	XG4	O6-C6-C5	-2.41	120.17	126.53
4	C	101	XG4	O1B-PB-N3A	-2.38	108.26	111.77
4	C	101	XG4	O2B-PB-O3B	2.34	112.45	104.64
4	B	1001	XG4	C4-C5-N7	-2.24	107.12	110.67
4	A	1001	XG4	C4-C5-N7	-2.24	107.12	110.67
4	B	1002	XG4	C4-C5-N7	-2.22	107.15	110.67
4	C	101	XG4	O3B-PG-O1G	-2.19	99.52	111.04
4	C	101	XG4	O6-C6-C5	-2.17	120.80	126.53
4	C	101	XG4	O5'-PA-O1A	-2.17	107.84	114.27
4	A	1001	XG4	C2'-C3'-C4'	2.13	107.11	102.80
4	B	1001	XG4	O1A-PA-N3A	-2.12	108.64	111.77
4	B	1002	XG4	O2A-PA-O5'	2.00	112.01	106.69
4	B	1001	XG4	C5-C6-N1	2.00	118.35	113.25

There are no chirality outliers.

All (36) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1001	XG4	PA-N3A-PB-O1B
4	A	1001	XG4	PB-O3B-PG-O3G
4	C	101	XG4	PB-N3A-PA-O1A
4	C	101	XG4	PA-N3A-PB-O1B
4	C	101	XG4	PG-O3B-PB-O2B
4	C	101	XG4	PB-O3B-PG-O2G
4	B	1001	XG4	PB-N3A-PA-O1A
4	B	1001	XG4	C5'-O5'-PA-O1A
4	B	1001	XG4	C5'-O5'-PA-O2A
4	B	1001	XG4	PA-N3A-PB-O1B
4	B	1001	XG4	PG-O3B-PB-O1B
4	B	1002	XG4	PA-N3A-PB-O1B
4	B	1002	XG4	PG-O3B-PB-O2B
4	B	1002	XG4	PB-O3B-PG-O2G
7	A	1009	GOL	O1-C1-C2-C3
4	A	1001	XG4	C3'-C4'-C5'-O5'

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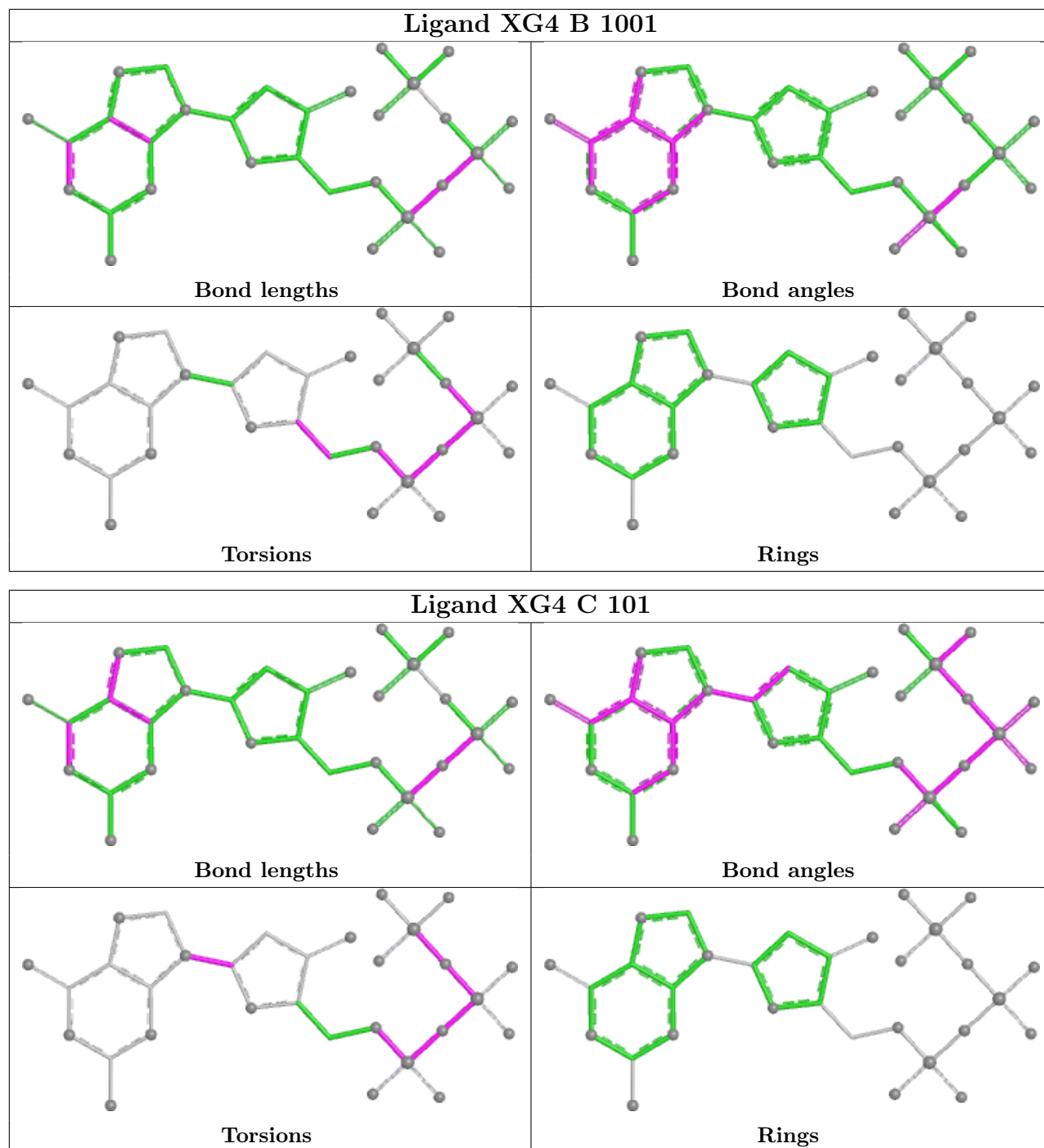
Mol	Chain	Res	Type	Atoms
4	B	1001	XG4	O4'-C4'-C5'-O5'
7	A	1008	GOL	O1-C1-C2-C3
7	A	1009	GOL	O1-C1-C2-O2
4	A	1001	XG4	O4'-C4'-C5'-O5'
4	B	1001	XG4	C3'-C4'-C5'-O5'
7	A	1008	GOL	O1-C1-C2-O2
4	C	101	XG4	O4'-C1'-N9-C4
4	A	1001	XG4	PB-O3B-PG-O1G
4	C	101	XG4	PB-O3B-PG-O1G
4	B	1002	XG4	PB-O3B-PG-O1G
4	B	1002	XG4	PB-N3A-PA-O5'
4	C	101	XG4	C5'-O5'-PA-O1A
4	B	1001	XG4	C5'-O5'-PA-N3A
4	B	1002	XG4	PB-N3A-PA-O1A
4	B	1001	XG4	PG-O3B-PB-O2B
4	C	101	XG4	C5'-O5'-PA-O2A
4	C	101	XG4	PB-O3B-PG-O3G
4	B	1002	XG4	PB-O3B-PG-O3G
4	B	1002	XG4	PG-O3B-PB-O1B
4	B	1001	XG4	PB-N3A-PA-O5'

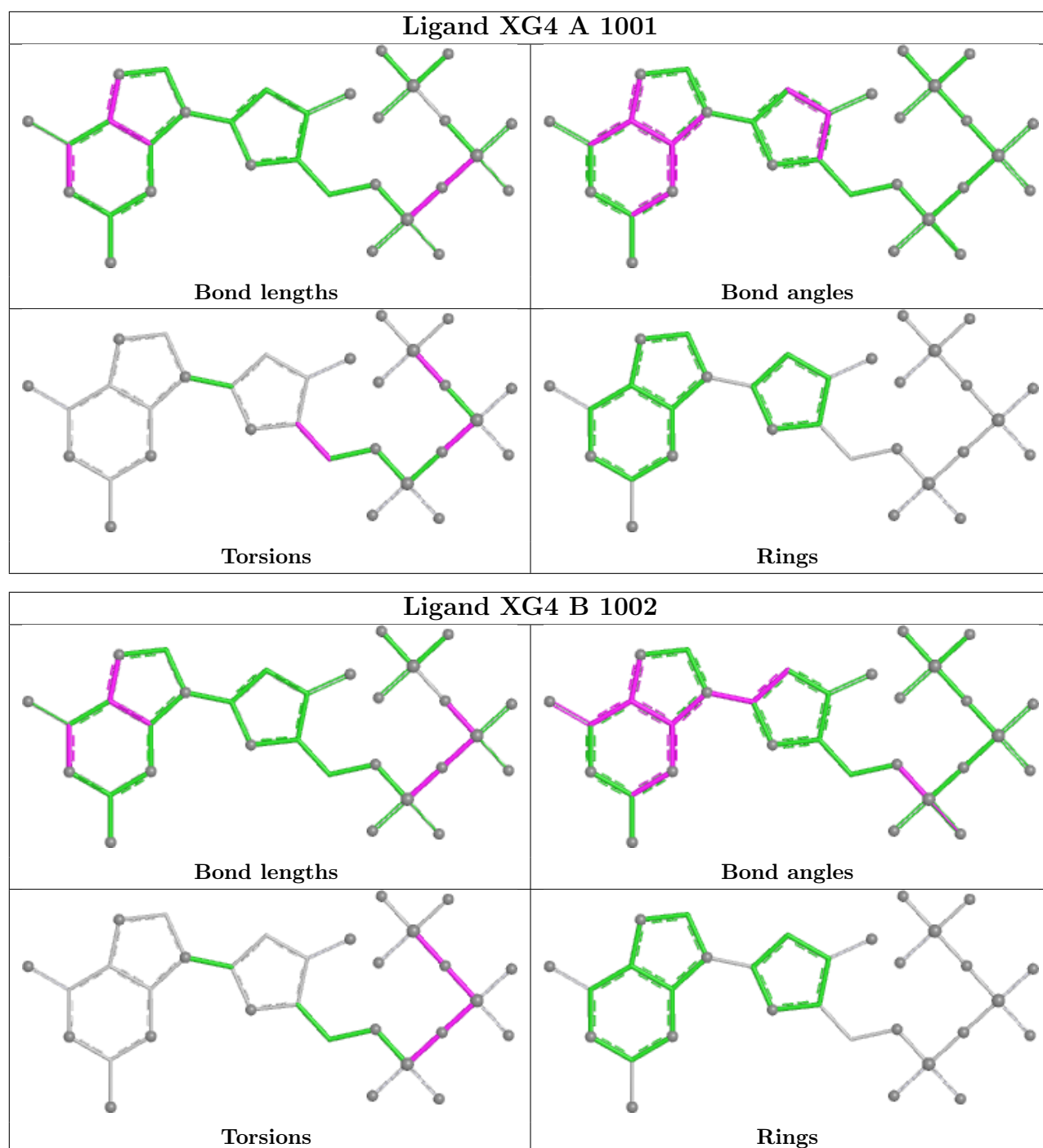
There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	101	XG4	2	0
4	B	1002	XG4	2	0
7	A	1008	GOL	4	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	901/903 (99%)	-0.46	7 (0%) 82 84	25, 46, 85, 118	2 (0%)
1	B	870/903 (96%)	0.11	16 (1%) 67 68	40, 81, 113, 135	1 (0%)
2	C	18/18 (100%)	-0.51	0 100 100	29, 44, 118, 128	0
2	E	18/18 (100%)	-0.45	0 100 100	50, 66, 118, 126	0
3	D	13/13 (100%)	-0.72	0 100 100	33, 42, 109, 113	0
3	F	13/13 (100%)	-0.48	0 100 100	55, 70, 111, 118	0
All	All	1833/1868 (98%)	-0.19	23 (1%) 75 75	25, 62, 109, 135	3 (0%)

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	257	TYR	4.7
1	A	901	PHE	3.7
1	B	303	LEU	3.6
1	A	256	MET	3.4
1	A	507	ASN	2.6
1	A	252	VAL	2.6
1	B	550	VAL	2.5
1	B	23	ASN	2.5
1	B	440[A]	HIS	2.5
1	B	820	ASP	2.4
1	B	46	ALA	2.4
1	B	535	ALA	2.3
1	B	473	THR	2.3
1	B	542	LEU	2.3
1	B	636	VAL	2.2
1	B	479	PHE	2.2
1	B	302	LYS	2.1
1	B	121	ASP	2.1
1	B	499	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	510	VAL	2.0
1	A	898	PHE	2.0
1	B	1	MET	2.0
1	B	22	SER	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](#)

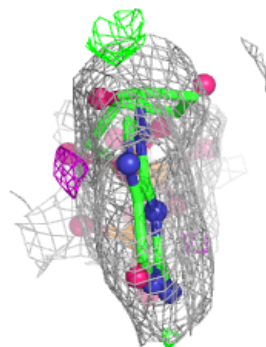
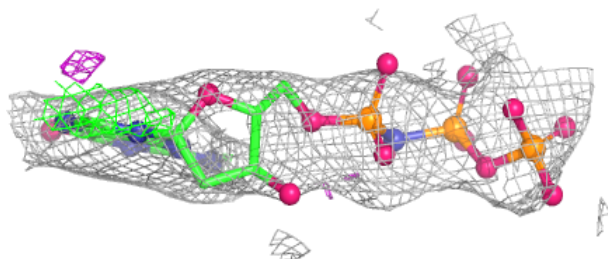
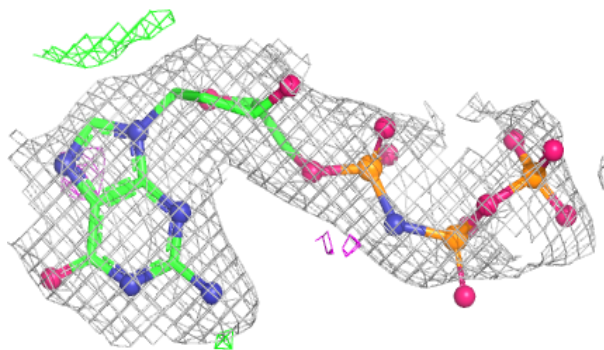
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
6	SO4	A	1005	5/5	0.78	0.09	114,116,118,118	0
4	XG4	B	1001	31/31	0.80	0.10	68,100,242,271	0
6	SO4	B	1003	5/5	0.81	0.19	93,95,98,98	0
7	GOL	A	1009	6/6	0.87	0.15	66,73,73,73	0
6	SO4	A	1006	5/5	0.88	0.10	109,110,111,113	0
6	SO4	A	1004	5/5	0.90	0.12	94,95,96,103	0
6	SO4	A	1007	5/5	0.91	0.12	136,137,137,139	0
4	XG4	B	1002	31/31	0.92	0.09	47,59,112,307	0
7	GOL	A	1008	6/6	0.93	0.13	44,50,62,69	0
4	XG4	A	1001	31/31	0.93	0.09	37,58,206,206	0
5	CA	B	1004	1/1	0.97	0.09	82,82,82,82	0
4	XG4	C	101	31/31	0.97	0.07	19,31,68,136	0
8	NA	A	1010	1/1	0.97	0.23	46,46,46,46	0
5	CA	A	1002	1/1	0.99	0.06	42,42,42,42	0
6	SO4	A	1003	5/5	0.99	0.04	47,47,48,53	0

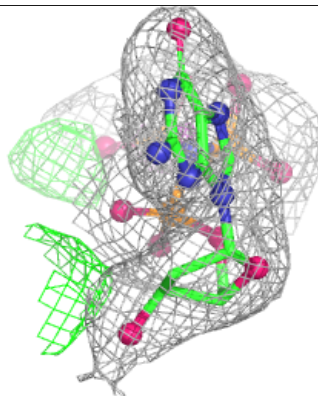
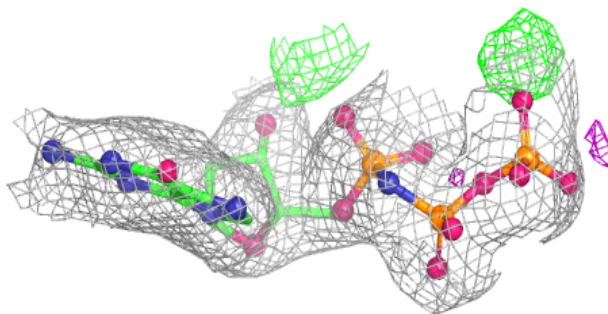
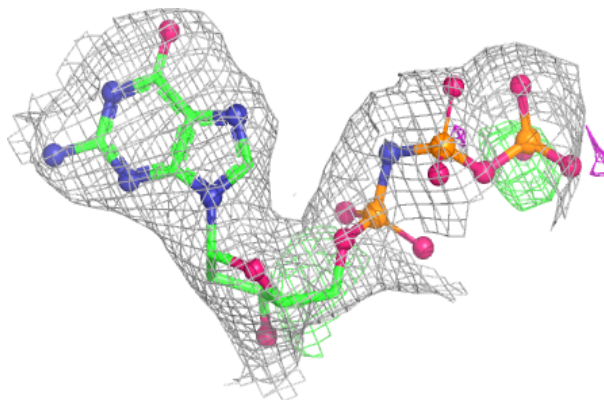
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around XG4 B 1001:

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

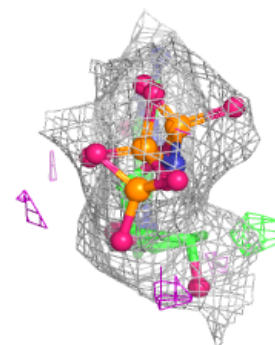
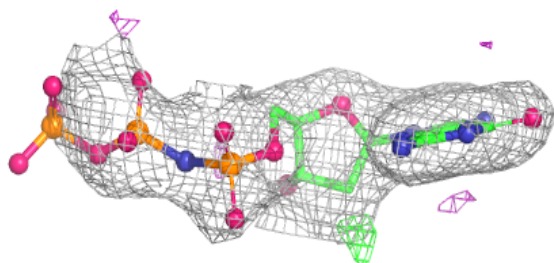
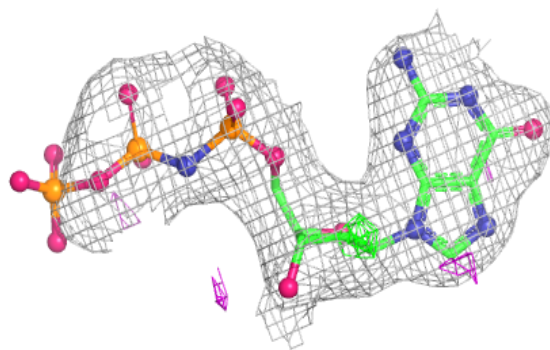
**Electron density around XG4 B 1002:**

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

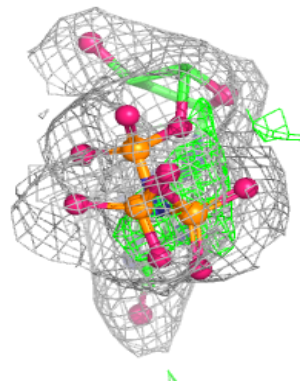
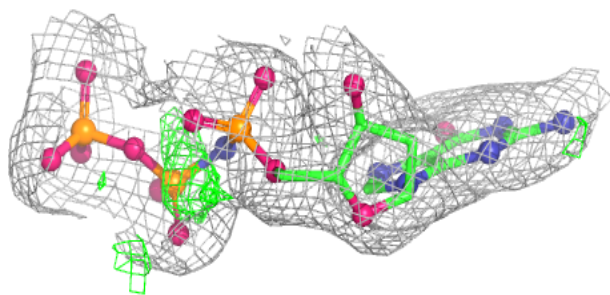
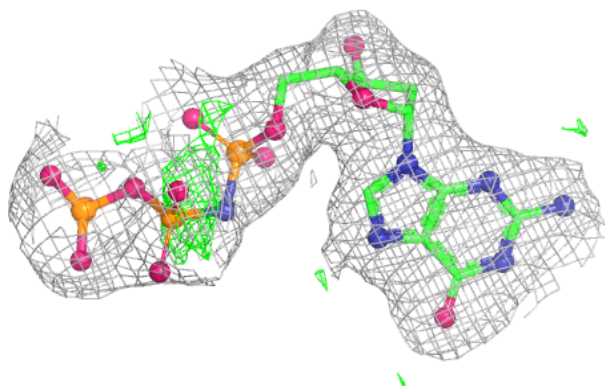


Electron density around XG4 A 1001:

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

**Electron density around XG4 C 101:**

$2mF_o-DF_c$ (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.