



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 18, 2022 – 12:15 PM JST

PDB ID : 7F4Y
Title : Crystal structure of replisomal dimer of DNA polymerase from bacteriophage RB69 with DNA duplexes
Authors : Youn, H.-S.; Park, J.; An, J.Y.; Lee, Y.; Eom, S.H.; Wang, J.
Deposited on : 2021-06-21
Resolution : 2.20 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.25
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.25

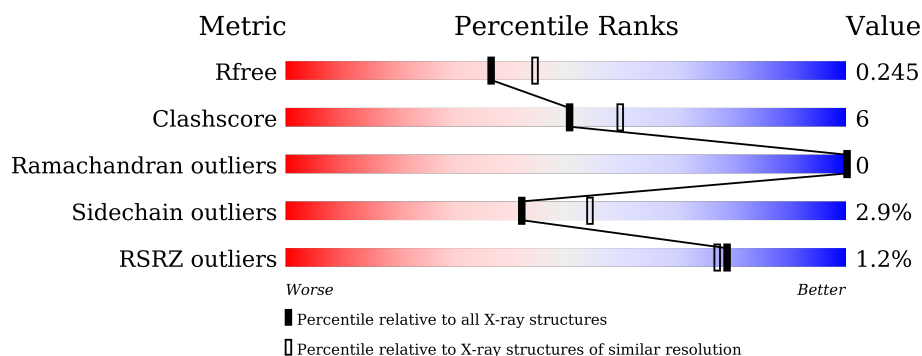
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	908	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="width: 100%; height: 10px; background-color: green;"></div> <div style="width: 100%; height: 10px; background-color: yellow;"></div> <div style="width: 100%; height: 10px; background-color: grey;"></div> </div> <div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="width: 100%; height: 10px; background-color: green;"></div> <div style="width: 100%; height: 10px; background-color: yellow;"></div> <div style="width: 100%; height: 10px; background-color: grey;"></div> </div> <div>87% 12% .</div>
1	B	908	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="width: 100%; height: 10px; background-color: green;"></div> <div style="width: 100%; height: 10px; background-color: yellow;"></div> <div style="width: 100%; height: 10px; background-color: grey;"></div> </div> <div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="width: 100%; height: 10px; background-color: green;"></div> <div style="width: 100%; height: 10px; background-color: yellow;"></div> <div style="width: 100%; height: 10px; background-color: grey;"></div> </div> <div>81% 18% ..</div>
2	S	19	<div> <div style="width: 100%; height: 10px; background-color: green;"></div> <div style="width: 100%; height: 10px; background-color: yellow;"></div> <div style="width: 100%; height: 10px; background-color: grey;"></div> </div> <div> <div style="width: 100%; height: 10px; background-color: green;"></div> <div style="width: 100%; height: 10px; background-color: yellow;"></div> <div style="width: 100%; height: 10px; background-color: grey;"></div> </div> <div>84% 16%</div>
2	T	19	<div> <div style="width: 100%; height: 10px; background-color: green;"></div> <div style="width: 100%; height: 10px; background-color: yellow;"></div> <div style="width: 100%; height: 10px; background-color: grey;"></div> </div> <div> <div style="width: 100%; height: 10px; background-color: green;"></div> <div style="width: 100%; height: 10px; background-color: yellow;"></div> <div style="width: 100%; height: 10px; background-color: grey;"></div> </div> <div>79% 16% 5%</div>
3	P	15	<div> <div style="width: 100%; height: 10px; background-color: green;"></div> <div style="width: 100%; height: 10px; background-color: yellow;"></div> <div style="width: 100%; height: 10px; background-color: grey;"></div> </div> <div> <div style="width: 100%; height: 10px; background-color: green;"></div> <div style="width: 100%; height: 10px; background-color: yellow;"></div> <div style="width: 100%; height: 10px; background-color: grey;"></div> </div> <div>67% 33%</div>
3	Q	15	<div> <div style="width: 100%; height: 10px; background-color: green;"></div> <div style="width: 100%; height: 10px; background-color: yellow;"></div> <div style="width: 100%; height: 10px; background-color: grey;"></div> </div> <div> <div style="width: 100%; height: 10px; background-color: green;"></div> <div style="width: 100%; height: 10px; background-color: yellow;"></div> <div style="width: 100%; height: 10px; background-color: grey;"></div> </div> <div>80% 20%</div>

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 18177 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	899	Total	C	N	O	S	0	13	0
			7411	4758	1241	1379	33			
1	B	901	Total	C	N	O	S	0	10	0
			7387	4745	1231	1379	32			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	expression tag	UNP Q38087
A	-3	ALA	-	expression tag	UNP Q38087
A	-2	MET	-	expression tag	UNP Q38087
A	-1	GLY	-	expression tag	UNP Q38087
A	0	SER	-	expression tag	UNP Q38087
A	222	ALA	ASP	engineered mutation	UNP Q38087
A	327	ALA	ASP	engineered mutation	UNP Q38087
B	-4	GLY	-	expression tag	UNP Q38087
B	-3	ALA	-	expression tag	UNP Q38087
B	-2	MET	-	expression tag	UNP Q38087
B	-1	GLY	-	expression tag	UNP Q38087
B	0	SER	-	expression tag	UNP Q38087
B	222	ALA	ASP	engineered mutation	UNP Q38087
B	327	ALA	ASP	engineered mutation	UNP Q38087

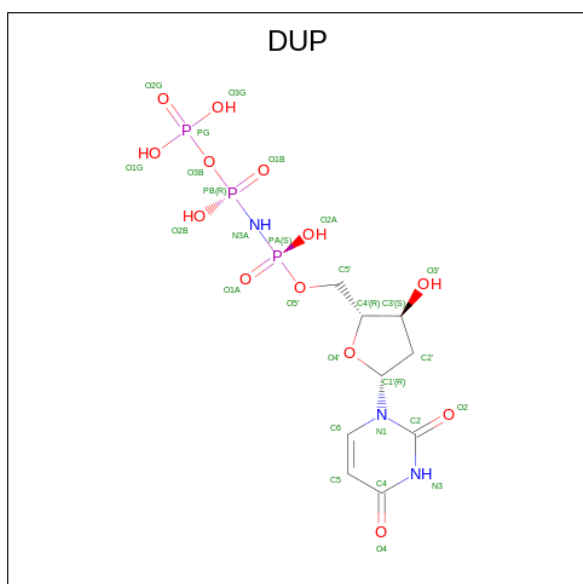
- Molecule 2 is a DNA chain called DNA (5'-D(*TP*CP*AP*AP*GP*TP*AP*AP*GP*CP*AP*GP*TP*CP*CP*GP*CP*TP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	T	18	Total	C	N	O	P	0	0	0
			367	174	69	106	18			
2	S	19	Total	C	N	O	P	0	0	0
			384	184	71	111	18			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	P	15	Total	C	N	O	P	0	0	0
			306	146	58	88	14			
3	Q	15	Total	C	N	O	P	0	0	0
			306	146	58	88	14			

- Molecule 4 is 2'-DEOXYURIDINE 5'-ALPHA,BETA-IMIDO-TRIPHOSPHATE (three-letter code: DUP) (formula: C₉H₁₆N₃O₁₃P₃).

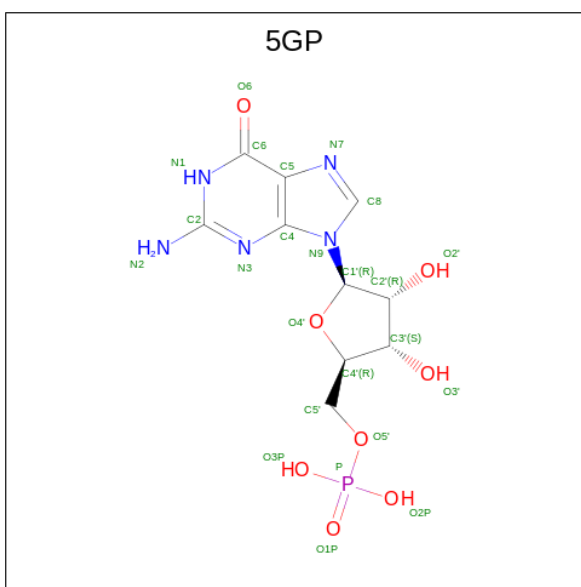


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			28	9	3	13	3		

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 6 is GUANOSINE-5'-MONOPHOSPHATE (three-letter code: 5GP) (formula: C₁₀H₁₄N₅O₈P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	P	0	0
			24	10	5	8	1		
6	B	1	Total	C	N	O	P	0	0
			24	10	5	8	1		

- Molecule 7 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	1	Total	Mg	0	0
			1	1		

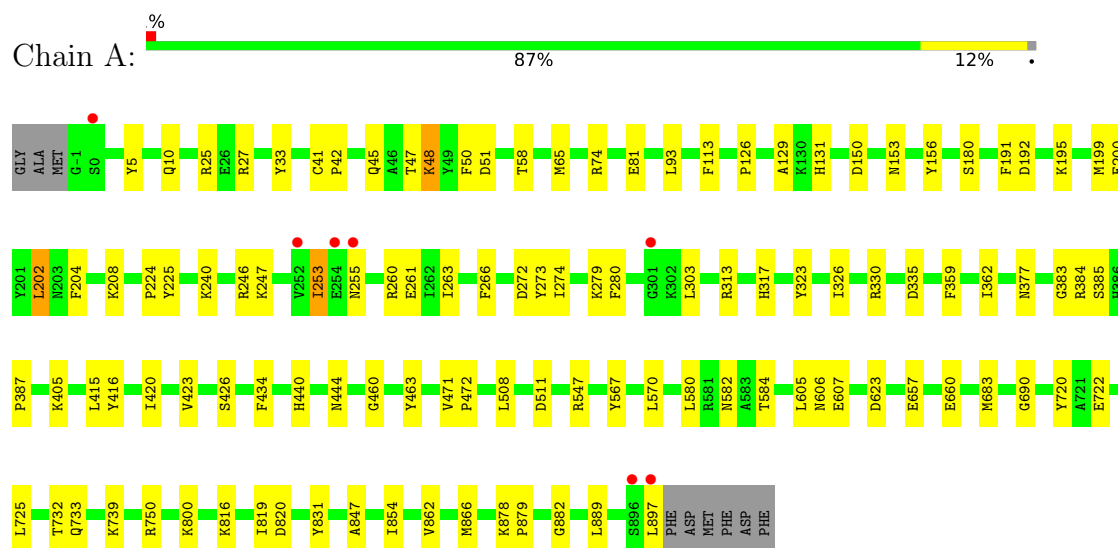
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	945	Total	O	0	3
			947	947		
8	T	76	Total	O	0	0
			76	76		
8	P	47	Total	O	0	0
			47	47		
8	B	723	Total	O	0	2
			725	725		
8	S	86	Total	O	0	1
			87	87		
8	Q	55	Total	O	0	0
			55	55		

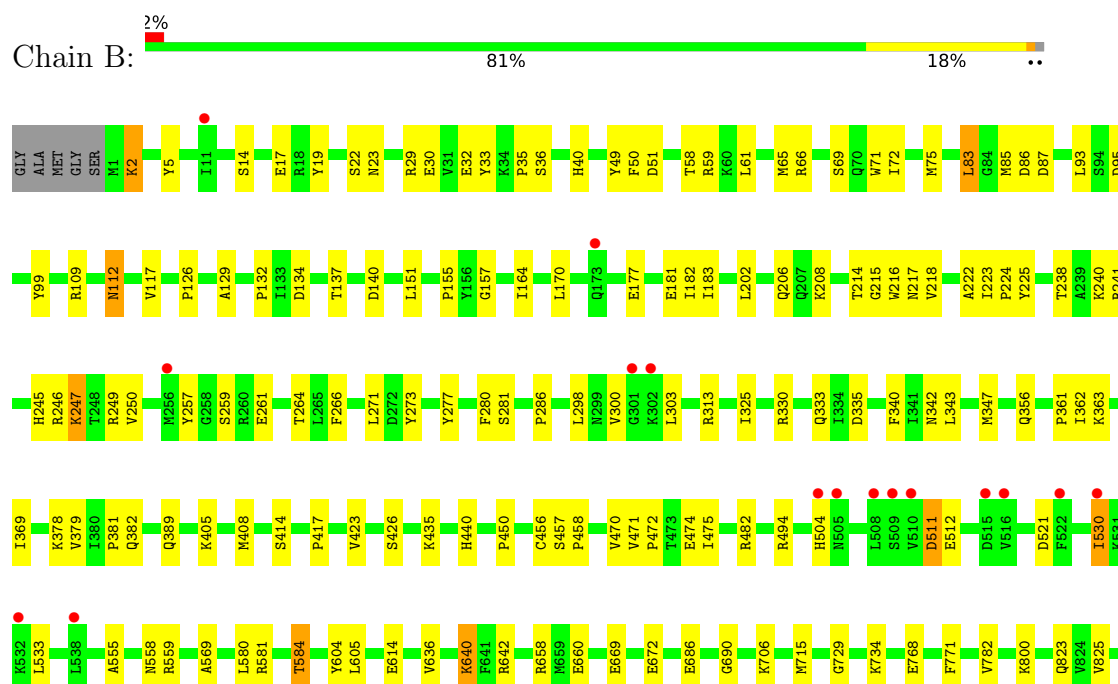
3 Residue-property plots [i](#)

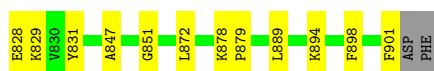
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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



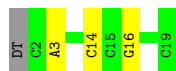
• Molecule 1: DNA polymerase





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

Chain T: 79% 16% 5%



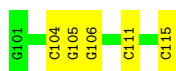
- Molecule 2: DNA (5'-D(*TP*CP*AP*AP*GP*TP*AP*AP*GP*CP*AP*GP*TP*CP*CP*GP*CP*TP*C)-3')

Chain S: 84% 16%



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

Chain P: 67% 33%



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

Chain Q: 80% 20%



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	164.44Å 164.44Å 165.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.01 – 2.20 45.13 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.9 (50.01-2.20) 99.9 (45.13-2.20)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.09 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.194 , 0.245 0.193 , 0.245	Depositor DCC
R_{free} test set	6484 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	38.8	Xtriage
Anisotropy	0.040	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 38.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.092 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	18177	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.99% 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: 5GP, MG, DUP, CA

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	0.62	0/7628	0.70	8/10303 (0.1%)
1	B	0.62	0/7598	0.68	0/10269
2	S	0.30	0/430	0.70	0/661
2	T	0.32	0/411	0.71	0/631
3	P	0.35	0/343	0.76	2/528 (0.4%)
3	Q	0.37	0/343	0.72	0/528
All	All	0.60	0/16753	0.69	10/22920 (0.0%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	607[A]	GLU	O-C-N	-6.55	112.22	122.70
1	A	607[B]	GLU	O-C-N	-6.55	112.22	122.70
1	A	606	ASN	O-C-N	5.83	132.02	122.70
3	P	115	DC	C1'-O4'-C4'	-5.70	104.40	110.10
1	A	260[A]	ARG	CA-C-O	5.68	132.03	120.10
1	A	260[B]	ARG	CA-C-O	5.68	132.03	120.10
1	A	260[A]	ARG	O-C-N	-5.48	113.92	122.70
1	A	260[B]	ARG	O-C-N	-5.48	113.92	122.70
3	P	111	DC	C1'-O4'-C4'	-5.17	104.93	110.10
1	A	383	GLY	C-N-CA	5.16	134.61	121.70

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	7411	0	7348	63	0
1	B	7387	0	7265	107	0
2	S	384	0	215	4	0
2	T	367	0	202	3	0
3	P	306	0	170	2	0
3	Q	306	0	170	2	0
4	A	28	0	12	4	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	24	0	12	0	0
6	B	24	0	12	2	0
7	B	1	0	0	0	0
8	A	947	0	0	9	0
8	B	725	0	0	14	0
8	P	47	0	0	1	0
8	Q	55	0	0	1	0
8	S	87	0	0	3	0
8	T	76	0	0	1	0
All	All	18177	0	15406	179	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 6.

All (179) 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:669:GLU:O	1:B:672[A]:GLU:HG3	1.57	1.04
4:A:1001:DUP:H5'1	4:A:1001:DUP:H6	1.40	1.04
1:A:81:GLU:HG3	1:A:384:ARG:NH2	1.82	0.95
1:B:828[B]:GLU:HG2	1:B:829:LYS:N	1.84	0.92
1:A:180:SER:HB2	8:A:1726:HOH:O	1.86	0.74
1:B:83:LEU:HD22	1:B:83:LEU:H	1.53	0.71
1:A:195:LYS:O	1:A:199[A]:MET:HG3	1.90	0.71
1:B:642:ARG:NH1	8:B:1106:HOH:O	2.24	0.70
1:B:71:TRP:O	1:B:75:MET:HG2	1.93	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:286:PRO:HB3	1:B:782:VAL:HG21	1.76	0.67
1:B:828[B]:GLU:HG2	1:B:829:LYS:H	1.59	0.65
1:B:5:TYR:CD2	1:B:93:LEU:HD22	2.31	0.64
1:B:273:TYR:OH	1:B:335:ASP:HA	1.98	0.64
1:B:471:VAL:HB	1:B:472:PRO:HD3	1.80	0.63
1:A:25[B]:ARG:NH2	8:A:1110:HOH:O	2.32	0.62
1:A:416:TYR:CD2	4:A:1001:DUP:H2'2	2.34	0.62
4:A:1001:DUP:H6	4:A:1001:DUP:C5'	2.23	0.62
1:B:17:GLU:OE1	1:B:19:TYR:HB3	2.00	0.62
1:B:414:SER:O	1:B:417:PRO:HD2	2.00	0.61
1:B:825:VAL:O	1:B:828[B]:GLU:HB3	1.99	0.61
1:A:279:LYS:NZ	8:A:1103:HOH:O	2.24	0.61
1:B:129:ALA:HA	1:B:225:TYR:CE2	2.36	0.60
1:A:204:PHE:CE2	1:A:208:LYS:HD3	2.35	0.60
1:B:555:ALA:O	1:B:559:ARG:HG2	2.02	0.60
1:B:472:PRO:O	1:B:475:ILE:HG22	2.02	0.60
1:B:280:PHE:HB3	1:B:340:PHE:CE1	2.37	0.59
1:A:279:LYS:HD3	1:A:280:PHE:CE1	2.37	0.59
1:A:303:LEU:HD22	1:A:323:TYR:CD1	2.37	0.59
1:A:153:ASN:HB2	1:A:192:ASP:O	2.03	0.58
1:B:686:GLU:O	1:B:715:MET:HA	2.03	0.58
1:B:423:VAL:O	1:B:423:VAL:HG12	2.03	0.58
1:B:878:LYS:HB3	1:B:879:PRO:HD3	1.84	0.57
2:S:15:DC:H2''	2:S:16:DG:C8	2.39	0.57
1:B:878:LYS:NZ	8:B:1109:HOH:O	2.28	0.57
1:B:83:LEU:HB3	1:B:379:VAL:HG12	1.87	0.56
1:A:195:LYS:NZ	8:A:1123:HOH:O	2.38	0.56
1:B:734:LYS:NZ	8:B:1119:HOH:O	2.33	0.56
1:A:5:TYR:CD2	1:A:93:LEU:HD22	2.40	0.56
1:B:95:ASP:OD2	6:B:1002:5GP:O3'	2.23	0.55
1:B:112:ASN:HA	1:B:214:THR:O	2.07	0.55
1:A:129:ALA:HA	1:A:225:TYR:CE2	2.43	0.54
1:B:330:ARG:HH11	1:B:333:GLN:HE22	1.55	0.54
1:B:580:LEU:O	1:B:584:THR:CG2	2.56	0.54
1:A:51:ASP:C	1:A:51:ASP:OD1	2.47	0.53
1:A:434:PHE:CZ	1:A:460:GLY:HA2	2.44	0.53
1:A:405:LYS:O	1:A:690:GLY:HA2	2.09	0.53
1:A:387:PRO:HB3	1:B:389:GLN:HB3	1.90	0.53
1:A:420:ILE:O	1:A:423:VAL:O	2.27	0.53
3:P:106:DG:N7	8:P:202:HOH:O	2.34	0.52
1:B:581:ARG:HD2	8:B:1173:HOH:O	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:T:16:DG:H5'	8:T:143:HOH:O	2.09	0.52
1:B:363:LYS:NZ	8:B:1107:HOH:O	2.25	0.52
1:A:273:TYR:OH	1:A:335:ASP:HA	2.10	0.52
1:A:580:LEU:O	1:A:584:THR:HG23	2.09	0.52
1:B:250:VAL:HG12	1:B:261:GLU:HG2	1.92	0.52
1:A:48:LYS:O	1:A:377:ASN:HB3	2.10	0.52
1:B:298:LEU:HB2	1:B:300:VAL:HG22	1.92	0.52
1:A:882:GLY:HA3	8:A:1464:HOH:O	2.09	0.52
1:B:51:ASP:OD1	1:B:51:ASP:C	2.48	0.52
1:A:359:PHE:HB3	2:T:3:DA:C4	2.44	0.51
1:A:47:THR:OG1	1:A:48:LYS:N	2.43	0.51
1:B:580:LEU:O	1:B:584:THR:HG22	2.10	0.51
1:B:706:LYS:O	1:B:729:GLY:HA3	2.10	0.51
1:B:218:VAL:HA	1:B:222:ALA:HB3	1.93	0.51
1:B:271:LEU:HD21	1:B:356:GLN:HA	1.92	0.51
1:B:470:VAL:O	1:B:474:GLU:HG2	2.11	0.51
1:B:823[A]:GLN:NE2	8:B:1113:HOH:O	2.31	0.51
1:B:208:LYS:HE3	8:B:1655:HOH:O	2.12	0.50
1:A:471:VAL:HB	1:A:472:PRO:HD3	1.92	0.50
1:B:50:PHE:O	1:B:378:LYS:HA	2.12	0.50
1:B:202:LEU:O	1:B:206:GLN:HG2	2.11	0.50
1:B:457:SER:OG	1:B:458:PRO:HD2	2.12	0.50
1:A:878:LYS:HB3	1:A:879:PRO:HD3	1.93	0.50
1:B:170:LEU:HA	1:B:177:GLU:HG3	1.94	0.49
3:P:104:DC:H2''	3:P:105:DG:C8	2.46	0.49
1:A:33:TYR:HA	8:A:1126:HOH:O	2.12	0.49
1:B:247:LYS:O	1:B:266:PHE:HB2	2.13	0.49
1:B:851:GLY:HA2	8:B:1174:HOH:O	2.11	0.49
1:B:343:LEU:HD12	1:B:558:ASN:HD22	1.78	0.48
1:B:35:PRO:HG3	1:B:65:MET:HA	1.95	0.48
1:A:434:PHE:CE1	1:A:460:GLY:HA2	2.49	0.48
1:A:326:ILE:O	1:A:330:ARG:HG2	2.13	0.48
1:B:69:SER:HA	1:B:72:ILE:HD12	1.94	0.48
1:B:800:LYS:HE2	8:S:154:HOH:O	2.13	0.48
1:A:126:PRO:HA	1:A:225:TYR:CD2	2.49	0.48
3:Q:106:DG:H4'	8:Q:229:HOH:O	2.13	0.48
2:S:16:DG:H5'	8:S:110:HOH:O	2.13	0.48
1:B:164:ILE:HD11	1:B:183:ILE:O	2.14	0.47
1:B:361:PRO:HG2	2:S:4:DA:O4'	2.14	0.47
1:A:313:ARG:HB3	8:A:1421:HOH:O	2.13	0.47
1:A:440:HIS:CE1	1:A:444:ASN:ND2	2.82	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:898:PHE:O	1:B:901:PHE:HB2	2.15	0.47
1:A:854:ILE:HD13	1:A:862:VAL:HG21	1.96	0.47
1:B:771:PHE:CG	1:B:872:LEU:HD13	2.49	0.47
1:B:828[B]:GLU:CG	1:B:829:LYS:H	2.25	0.47
1:B:240:LYS:O	1:B:246:ARG:HA	2.15	0.47
1:B:382:GLN:HA	8:B:1112:HOH:O	2.14	0.46
1:A:81:GLU:CG	1:A:384:ARG:NH2	2.68	0.46
1:A:41:CYS:HB2	1:A:42:PRO:HD2	1.97	0.46
1:B:2:LYS:HG3	1:B:99:TYR:CE2	2.50	0.46
1:B:604:TYR:OH	1:B:658[B]:ARG:HB3	2.15	0.46
1:B:347:MET:HB2	1:B:558:ASN:HD21	1.80	0.46
1:B:65:MET:N	8:B:1150:HOH:O	2.47	0.46
1:B:636:VAL:HG12	1:B:640:LYS:HE2	1.98	0.46
1:A:202:LEU:HD23	1:A:202:LEU:HA	1.82	0.46
1:A:224:PRO:HA	1:A:263:ILE:HG13	1.98	0.46
1:A:463:TYR:OH	1:A:582:ASN:ND2	2.49	0.46
1:B:126:PRO:HA	1:B:225:TYR:HD2	1.80	0.46
1:B:181:GLU:HG2	1:B:182:ILE:HG23	1.98	0.45
1:B:894:LYS:HE2	8:B:1122:HOH:O	2.16	0.45
1:B:157:GLY:HA3	1:B:313:ARG:HH12	1.82	0.45
1:B:215:GLY:HA3	1:B:218:VAL:CG1	2.47	0.45
1:B:245:HIS:C	1:B:247:LYS:N	2.68	0.45
1:B:440[A]:HIS:CE1	8:B:1244:HOH:O	2.69	0.45
1:A:416:TYR:CG	4:A:1001:DUP:H2'2	2.51	0.45
1:B:450:PRO:HB2	1:B:456:CYS:SG	2.57	0.45
1:B:33:TYR:HB3	1:B:65:MET:HE2	1.97	0.45
1:B:241[B]:ARG:HE	1:B:246:ARG:HD2	1.82	0.45
1:B:471:VAL:HB	1:B:472:PRO:CD	2.46	0.45
1:A:253:ILE:HD13	1:A:253:ILE:HA	1.69	0.45
1:A:725:LEU:HD11	1:A:750:ARG:HB2	1.98	0.45
1:B:69:SER:O	1:B:72:ILE:HB	2.17	0.45
1:B:216:TRP:O	1:B:217:ASN:HB2	2.16	0.45
1:A:831:TYR:O	1:A:847:ALA:HA	2.17	0.44
1:B:405:LYS:O	1:B:690:GLY:HA2	2.17	0.44
1:B:604:TYR:OH	1:B:658[A]:ARG:HB3	2.16	0.44
1:A:272:ASP:OD1	1:A:274:ILE:HG22	2.17	0.44
1:A:732:THR:OG1	1:A:733:GLN:NE2	2.31	0.44
1:B:277:TYR:O	1:B:281:SER:HB3	2.18	0.44
1:A:800:LYS:HE3	2:T:14:DC:H4'	1.99	0.43
1:A:126:PRO:HA	1:A:225:TYR:HD2	1.83	0.43
1:A:191:PHE:HZ	1:A:200:GLU:HG2	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:109:ARG:HH11	1:B:140:ASP:CG	2.20	0.43
1:B:132:PRO:HG2	1:B:155:PRO:HD2	1.99	0.43
3:Q:108:DC:H2''	3:Q:109:DT:H5'	1.99	0.43
1:A:862:VAL:O	1:A:866:MET:HG3	2.19	0.43
1:B:249:ARG:HG2	1:B:264:THR:HB	2.01	0.43
1:B:271:LEU:CD2	1:B:356:GLN:HA	2.49	0.43
1:B:134:ASP:HA	1:B:151:LEU:HB3	1.99	0.43
1:B:223:ILE:N	1:B:224:PRO:HD2	2.33	0.43
1:B:482:ARG:HG3	1:B:559:ARG:HB2	2.01	0.43
1:B:831:TYR:O	1:B:847:ALA:HA	2.18	0.43
1:A:415:LEU:HD22	1:A:623:ASP:HB3	2.01	0.42
1:A:567:TYR:O	1:A:570:LEU:HB2	2.20	0.42
1:B:238:THR:O	1:B:241[A]:ARG:HB2	2.18	0.42
1:A:247:LYS:HE3	1:A:266:PHE:CZ	2.54	0.42
1:A:660:GLU:OE2	1:A:683:MET:O	2.37	0.42
1:A:240:LYS:O	1:A:246:ARG:HA	2.19	0.42
1:A:720:TYR:HB3	1:A:722:GLU:O	2.19	0.42
1:B:19:TYR:CZ	1:B:29:ARG:NH1	2.87	0.42
1:B:40:HIS:HB3	8:B:1618:HOH:O	2.20	0.42
1:B:245:HIS:C	1:B:247:LYS:H	2.22	0.42
2:S:16:DG:N7	8:S:102:HOH:O	2.36	0.42
1:B:605:LEU:HD23	1:B:605:LEU:HA	1.89	0.42
1:B:14:SER:HB3	1:B:30:GLU:HG3	2.02	0.42
1:B:494:ARG:NH1	1:B:521:ASP:OD1	2.50	0.42
1:B:238:THR:O	1:B:241[B]:ARG:HB2	2.19	0.41
1:B:423:VAL:O	1:B:423:VAL:CG1	2.67	0.41
1:B:530:ILE:HA	1:B:533:LEU:HD12	2.02	0.41
1:B:85:MET:HG3	1:B:87:ASP:O	2.19	0.41
1:B:381:PRO:CB	8:B:1429:HOH:O	2.68	0.41
1:A:10:GLN:HG3	1:A:65:MET:SD	2.61	0.41
1:A:816:LYS:NZ	8:A:1168:HOH:O	2.52	0.41
1:B:362:ILE:HD11	1:B:569:ALA:HA	2.01	0.41
1:B:303:LEU:H	1:B:303:LEU:HD12	1.86	0.41
1:A:74:ARG:HD2	8:A:1627:HOH:O	2.21	0.41
1:B:32:GLU:CG	1:B:32:GLU:O	2.68	0.41
1:A:47:THR:O	1:A:50:PHE:CE1	2.74	0.41
1:A:547:ARG:HD3	1:A:547:ARG:HA	1.74	0.41
1:A:657:GLU:OE2	1:A:657:GLU:HA	2.21	0.41
1:B:35:PRO:O	1:B:61:LEU:HD12	2.20	0.41
1:B:511:ASP:HB2	1:B:533:LEU:HD23	2.03	0.41
1:B:117:VAL:HG11	1:B:225:TYR:CZ	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:150:ASP:OD2	1:A:317:HIS:CE1	2.74	0.40
1:A:605:LEU:HD23	1:A:605:LEU:HA	1.87	0.40
1:B:49:TYR:OH	6:B:1002:5GP:O3P	2.28	0.40
1:A:131:HIS:HD2	1:A:156:TYR:OH	2.04	0.40
1:B:369:ILE:HG12	1:B:474:GLU:HG3	2.04	0.40
1:B:137:THR:HG21	1:B:325:ILE:HA	2.02	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	910/908 (100%)	876 (96%)	34 (4%)	0	100	100
1	B	909/908 (100%)	855 (94%)	54 (6%)	0	100	100
All	All	1819/1816 (100%)	1731 (95%)	88 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	806/802 (100%)	787 (98%)	19 (2%)	49	62
1	B	798/802 (100%)	771 (97%)	27 (3%)	37	47

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1604/1604 (100%)	1558 (97%)	46 (3%)	42 54

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	ARG
1	A	45	GLN
1	A	48	LYS
1	A	58	THR
1	A	113	PHE
1	A	202	LEU
1	A	253	ILE
1	A	255	ASN
1	A	261	GLU
1	A	362	ILE
1	A	385	SER
1	A	426	SER
1	A	508	LEU
1	A	511	ASP
1	A	739	LYS
1	A	819	ILE
1	A	820	ASP
1	A	889	LEU
1	A	897	LEU
1	B	2	LYS
1	B	22	SER
1	B	23	ASN
1	B	36	SER
1	B	58	THR
1	B	59	ARG
1	B	66	ARG
1	B	83	LEU
1	B	86	ASP
1	B	112	ASN
1	B	247	LYS
1	B	257	TYR
1	B	259	SER
1	B	342	ASN
1	B	408	MET
1	B	426	SER
1	B	435	LYS
1	B	504	HIS

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Mol	Chain	Res	Type
1	B	511	ASP
1	B	512	GLU
1	B	530	ILE
1	B	584	THR
1	B	614	GLU
1	B	640	LYS
1	B	660	GLU
1	B	768	GLU
1	B	889	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	131	HIS
1	A	206	GLN
1	A	339	GLN
1	A	440	HIS
1	A	444	ASN
1	A	507	ASN
1	A	546	GLN
1	A	582	ASN
1	A	733	GLN
1	A	761	GLN
1	B	128	GLN
1	B	131	HIS
1	B	333	GLN
1	B	354	GLN
1	B	444	ASN
1	B	546	GLN
1	B	582	ASN
1	B	678	GLN
1	B	761	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 3 are monoatomic - leaving 3 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	DUP	A	1001	5	28,29,29	2.07	8 (28%)	37,45,45	1.55	7 (18%)
6	5GP	A	1003	-	22,26,26	1.04	2 (9%)	27,40,40	2.33	4 (14%)
6	5GP	B	1002	-	22,26,26	1.06	1 (4%)	27,40,40	2.26	4 (14%)

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	DUP	A	1001	5	-	5/19/34/34	0/2/2/2
6	5GP	A	1003	-	-	2/6/26/26	0/3/3/3
6	5GP	B	1002	-	-	1/6/26/26	0/3/3/3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1001	DUP	C6-C5	-6.90	1.34	1.52
4	A	1001	DUP	C6-N1	-5.12	1.37	1.47
6	B	1002	5GP	C6-N1	3.94	1.39	1.33
6	A	1003	5GP	C6-N1	3.67	1.39	1.33
4	A	1001	DUP	C5-C4	-2.90	1.43	1.50
4	A	1001	DUP	PA-O2A	-2.69	1.49	1.56
4	A	1001	DUP	C1'-N1	2.63	1.49	1.45
4	A	1001	DUP	PB-O2B	-2.33	1.50	1.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1001	DUP	PB-O1B	2.30	1.49	1.46
4	A	1001	DUP	PA-O1A	2.19	1.49	1.46
6	A	1003	5GP	C8-N7	-2.06	1.31	1.34

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	1003	5GP	C5-C6-N1	-8.93	111.22	123.43
6	B	1002	5GP	C5-C6-N1	-8.89	111.28	123.43
6	A	1003	5GP	C2-N1-C6	6.04	125.52	115.93
6	B	1002	5GP	C2-N1-C6	5.91	125.31	115.93
4	A	1001	DUP	O2B-PB-O1B	4.21	118.75	109.92
4	A	1001	DUP	C5-C6-N1	3.66	123.66	111.61
4	A	1001	DUP	O2A-PA-O1A	3.60	117.47	109.92
6	A	1003	5GP	C2-N3-C4	-2.91	112.03	115.36
6	B	1002	5GP	C2-N3-C4	-2.82	112.14	115.36
6	A	1003	5GP	N3-C2-N1	-2.73	123.58	127.22
6	B	1002	5GP	N3-C2-N1	-2.69	123.63	127.22
4	A	1001	DUP	C4-N3-C2	2.45	127.82	125.79
4	A	1001	DUP	O5'-PA-O1A	-2.10	106.16	114.24
4	A	1001	DUP	O3B-PB-N3A	-2.08	100.82	106.59
4	A	1001	DUP	C5-C4-N3	-2.01	114.39	116.65

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1001	DUP	C3'-C4'-C5'-O5'
4	A	1001	DUP	PA-N3A-PB-O1B
4	A	1001	DUP	PB-O3B-PG-O3G
6	A	1003	5GP	O4'-C4'-C5'-O5'
6	A	1003	5GP	C3'-C4'-C5'-O5'
4	A	1001	DUP	O4'-C4'-C5'-O5'
6	B	1002	5GP	O4'-C4'-C5'-O5'
4	A	1001	DUP	PB-O3B-PG-O2G

There are no ring outliers.

2 monomers are involved in 6 short contacts:

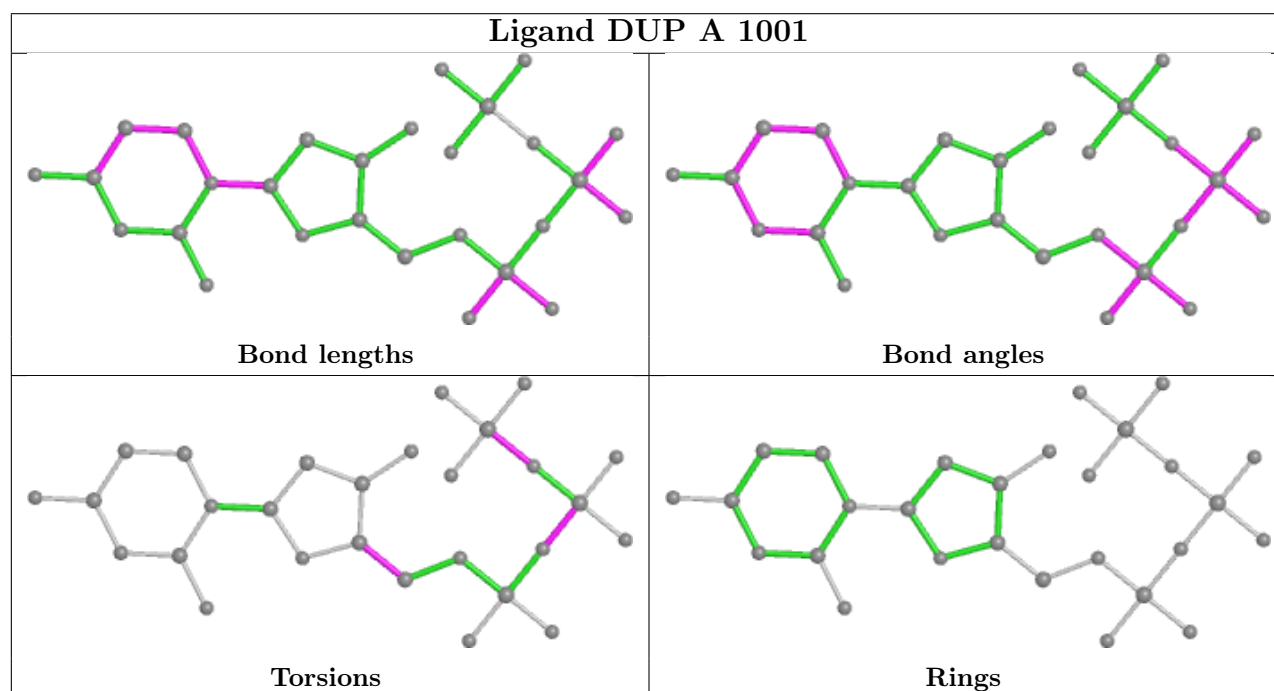
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1001	DUP	4	0

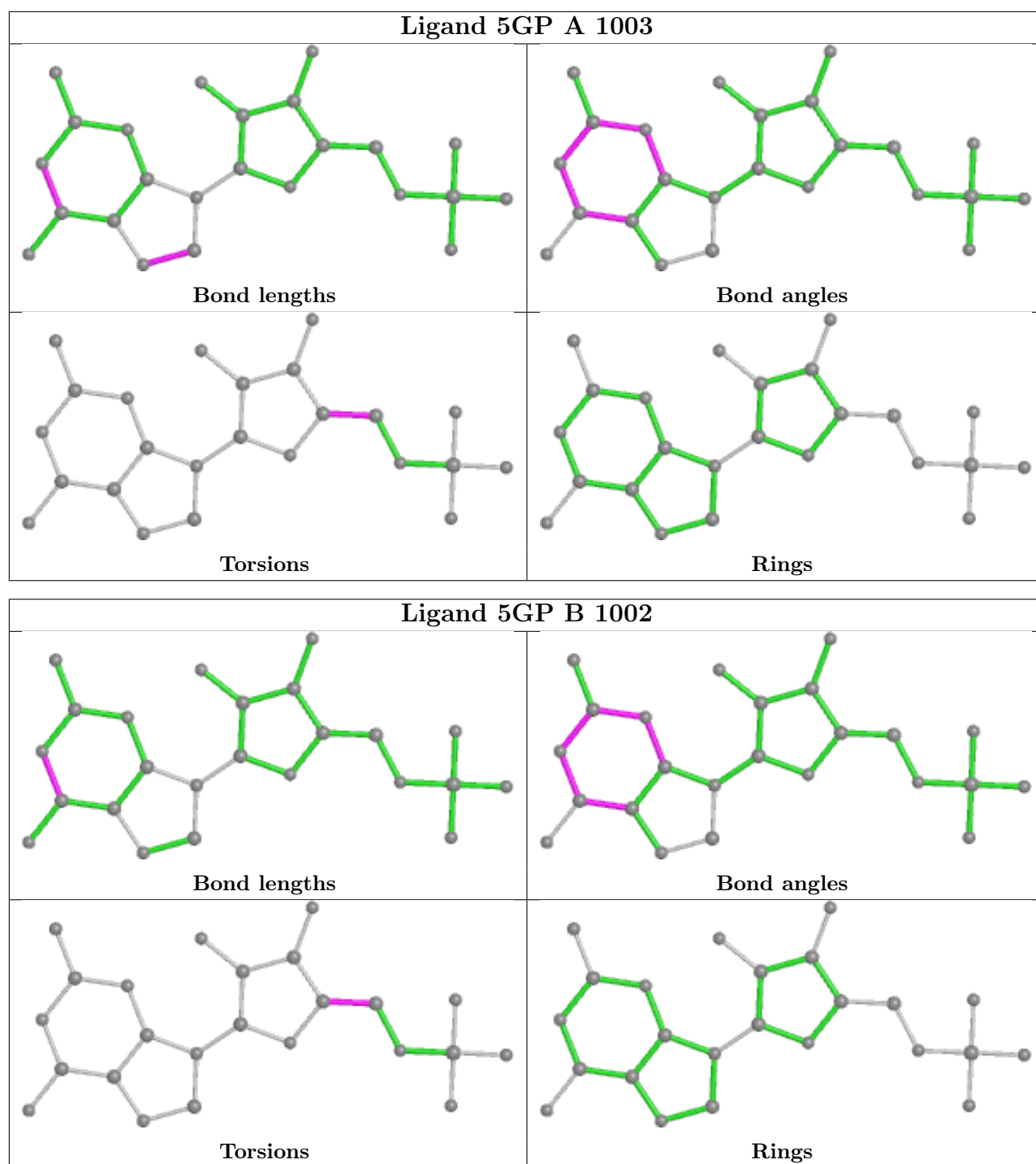
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	1002	5GP	2	0

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	899/908 (99%)	-0.72	7 (0%) 86 85	23, 42, 92, 145	0
1	B	901/908 (99%)	-0.42	16 (1%) 68 66	24, 52, 112, 145	0
2	S	19/19 (100%)	-1.06	0 100 100	29, 36, 67, 82	0
2	T	18/19 (94%)	-0.94	0 100 100	26, 33, 80, 103	0
3	P	15/15 (100%)	-1.08	0 100 100	24, 36, 61, 63	0
3	Q	15/15 (100%)	-1.09	0 100 100	27, 37, 62, 69	0
All	All	1867/1884 (99%)	-0.59	23 (1%) 79 77	23, 46, 103, 145	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	510	VAL	7.2
1	B	301	GLY	5.6
1	B	504	HIS	5.1
1	B	532	LYS	3.9
1	A	254	GLU	3.8
1	B	256	MET	3.7
1	A	0	SER	3.5
1	B	509	SER	3.4
1	B	516	VAL	3.2
1	B	538	LEU	3.0
1	A	255	ASN	2.8
1	B	508	LEU	2.7
1	A	252	VAL	2.6
1	B	11	ILE	2.4
1	B	505	ASN	2.4
1	B	173	GLN	2.4
1	B	530	ILE	2.3
1	A	897	LEU	2.3
1	A	896	SER	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	302	LYS	2.1
1	A	301	GLY	2.0
1	B	522	PHE	2.0
1	B	515	ASP	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 monosaccharides in this entry.

6.4 Ligands [i](#)

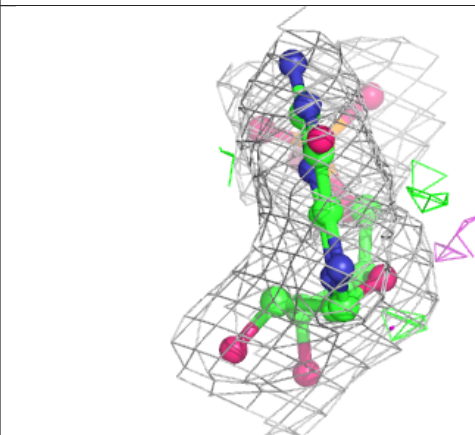
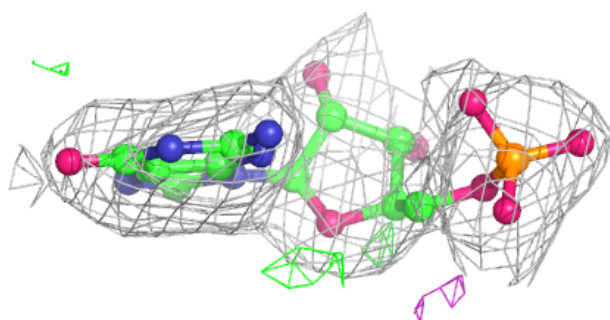
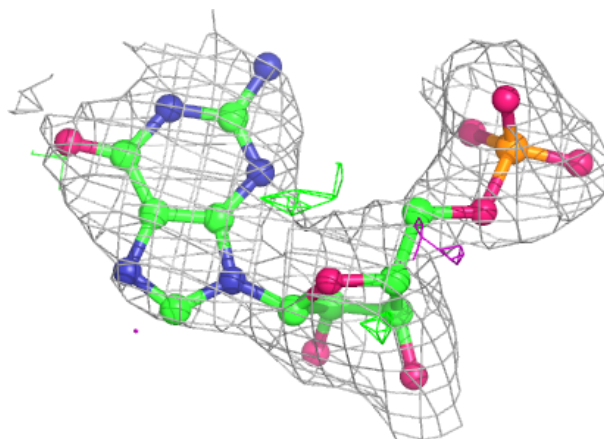
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	CA	B	1001	1/1	0.89	0.18	108,108,108,108	0
5	CA	A	1002	1/1	0.90	0.07	63,63,63,63	0
6	5GP	B	1002	24/24	0.91	0.12	59,75,90,102	0
6	5GP	A	1003	24/24	0.94	0.10	46,54,73,81	0
4	DUP	A	1001	28/28	0.96	0.09	28,38,50,52	0
7	MG	B	1003	1/1	0.97	0.17	64,64,64,64	0

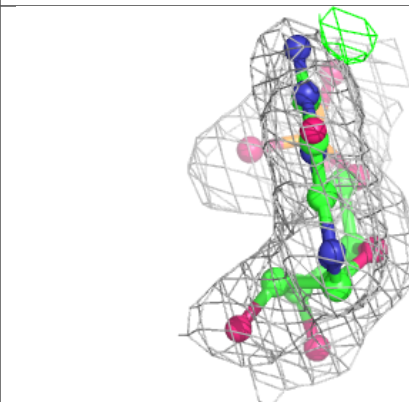
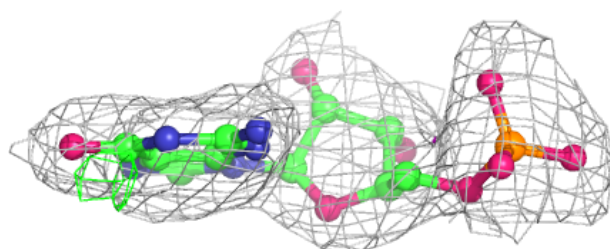
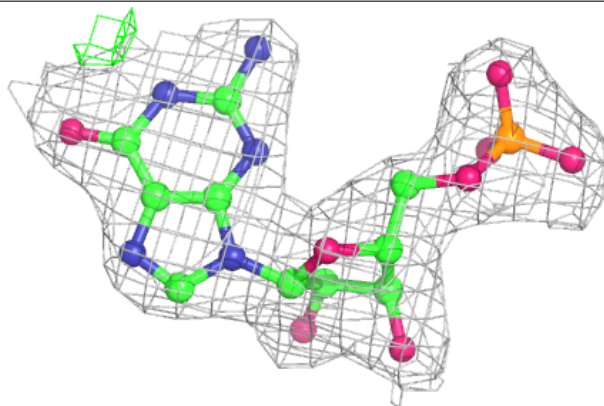
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 5GP 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)

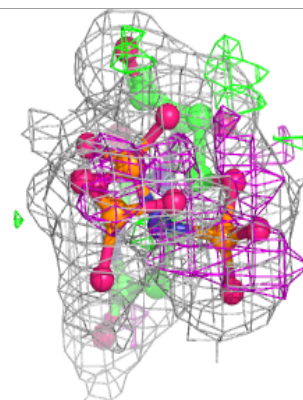
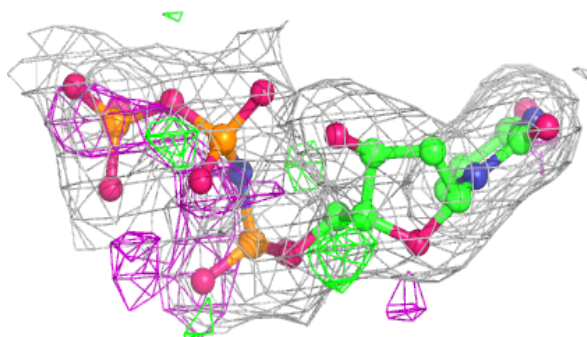
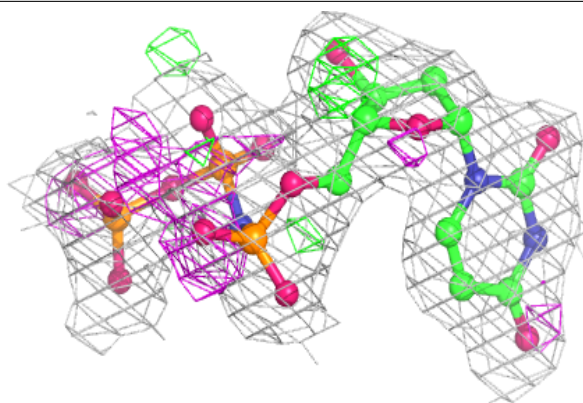
**Electron density around 5GP A 1003:**

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



Electron density around DUP 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.