



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 07:53 am BST

PDB ID : 5ZLV
Title : DNA polymerase IV - DNA ternary complex with 50mM MgCl₂
Authors : Kottur, J.; Nair, D.T.
Deposited on : 2018-03-29
Resolution : 2.35 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 : 2.11
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) : 2.11

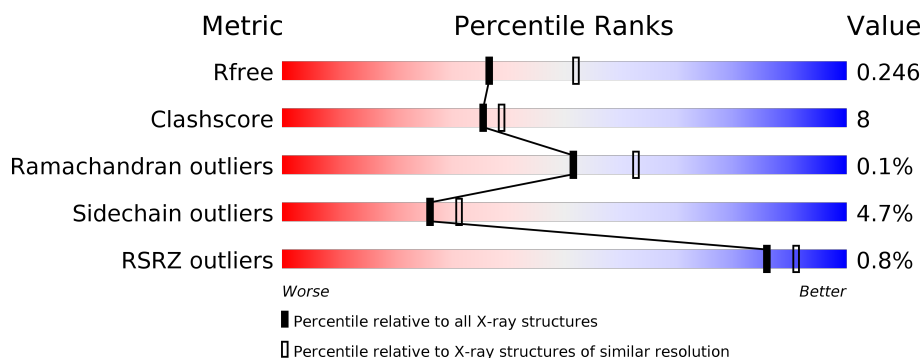
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

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 on 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	352	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 16%, green 79%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 1% 79% 16% </div> </div>
1	F	352	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, yellow 17%, green 79%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 17% 79% </div> </div>
2	B	18	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 6%, orange 1%, yellow 22%, green 72%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 6% 72% 22% </div> </div>
2	G	18	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 6%, orange 1%, yellow 22%, green 67%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 6% 67% 22% </div> </div>
3	C	19	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, yellow 42%, green 47%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 42% 47% </div> </div>
3	H	19	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, yellow 21%, green 53%, grey 26%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 21% 53% 26% </div> </div>

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
5	PO4	F	903	-	-	X	-
6	DPO	A	903	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7158 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 IV.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	F	342	Total	C	N	O	S	0	0	0
			2691	1698	495	484	14			
1	A	342	Total	C	N	O	S	0	0	0
			2691	1698	495	484	14			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	0	GLY	-	expression tag	UNP Q47155
F	1	SER	-	expression tag	UNP Q47155
A	0	GLY	-	expression tag	UNP Q47155
A	1	SER	-	expression tag	UNP Q47155

- Molecule 2 is a DNA chain called DTN2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	16	Total	C	N	O	P	0	0	0
			328	155	61	96	16			
2	B	17	Total	C	N	O	P	0	0	0
			347	164	64	102	17			

- Molecule 3 is a DNA chain called DTN1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	14	Total	C	N	O	P	0	0	0
			285	135	51	85	14			
3	C	17	Total	C	N	O	P	0	0	0
			348	165	63	103	17			

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

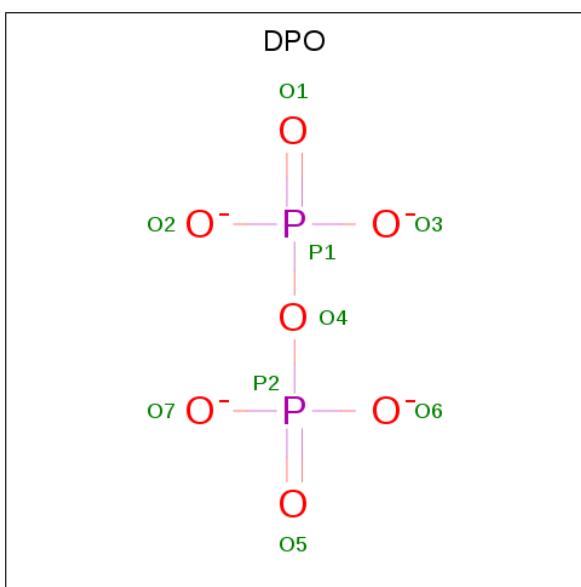
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	1	Total	Mg	0	0
			1	1		
4	A	2	Total	Mg	0	0
			2	2		
4	F	2	Total	Mg	0	0
			2	2		

- Molecule 5 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	F	1	Total	O	P	0	0
			5	4	1		
5	F	1	Total	O	P	0	0
			5	4	1		

- Molecule 6 is DIPHOSPHATE (three-letter code: DPO) (formula: O₇P₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	P	0	0
			9	7	2		

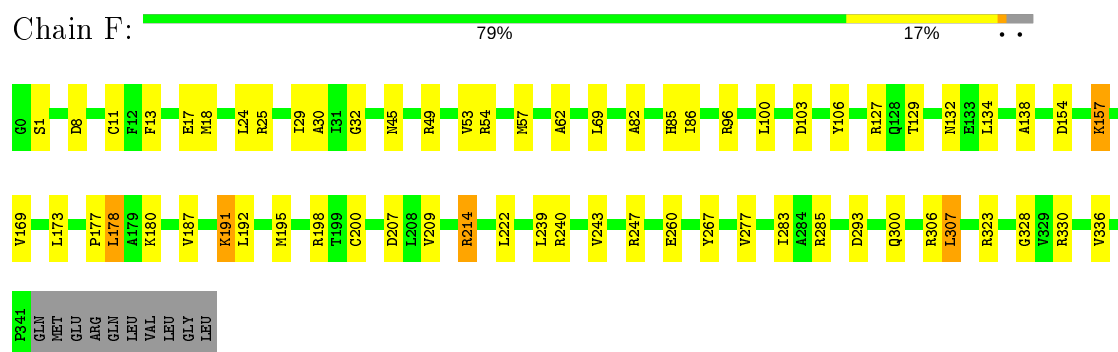
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	F	166	Total	O	0	0
			166	166		
7	A	159	Total	O	0	0
			159	159		
7	G	29	Total	O	0	0
			29	29		
7	H	22	Total	O	0	0
			22	22		
7	B	28	Total	O	0	0
			28	28		
7	C	40	Total	O	0	0
			40	40		

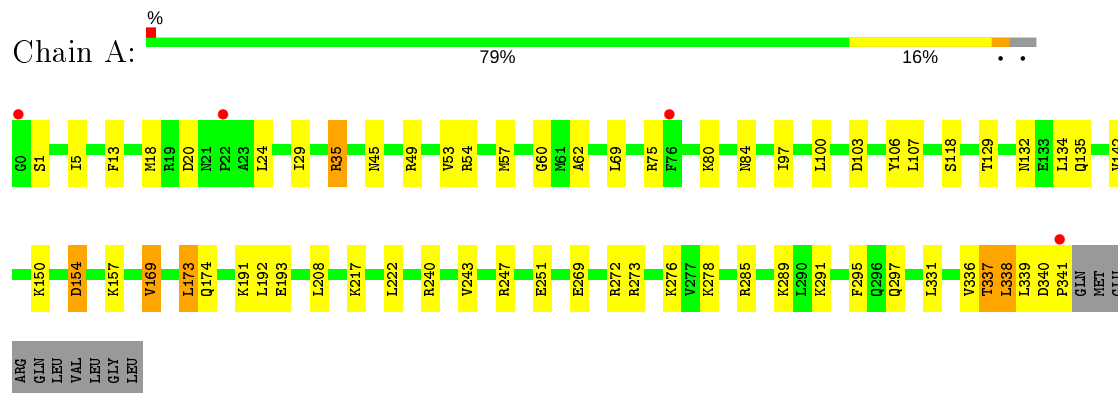
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

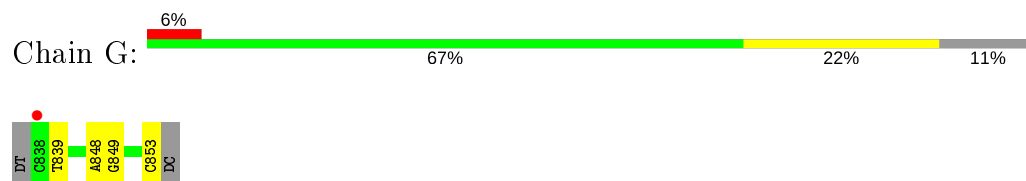
• Molecule 1: DNA polymerase IV



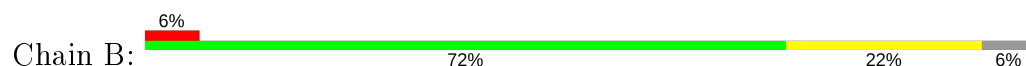
• Molecule 1: DNA polymerase IV



• Molecule 2: DTN2

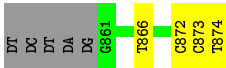


• Molecule 2: DTN2





● Molecule 3: DTN1



● Molecule 3: DTN1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	86.31Å 57.24Å 110.39Å 90.00° 94.74° 90.00°	Depositor
Resolution (Å)	50.78 – 2.35 65.20 – 2.35	Depositor EDS
% Data completeness (in resolution range)	96.7 (50.78-2.35) 98.3 (65.20-2.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.97 (at 2.34Å)	Xtriage
Refinement program	PHENIX (1.10.1 _2155: ???)	Depositor
R, R_{free}	0.188 , 0.248 0.186 , 0.246	Depositor DCC
R_{free} test set	2253 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	32.1	Xtriage
Anisotropy	0.609	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 42.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7158	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.96% 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: DPO, PO4, MG

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	0.37	0/2742	0.56	0/3703
1	F	0.42	0/2742	0.61	0/3703
2	B	0.80	0/388	0.90	0/596
2	G	0.97	1/367 (0.3%)	0.93	0/564
3	C	0.94	1/389 (0.3%)	0.96	0/598
3	H	0.93	0/318	0.98	0/488
All	All	0.54	2/6946 (0.0%)	0.69	0/9652

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	839	DT	C3'-O3'	-5.98	1.36	1.44
3	C	867	DA	C3'-O3'	-5.83	1.36	1.44

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2691	0	2750	42	1
1	F	2691	0	2750	49	0
2	B	347	0	191	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	G	328	0	180	3	0
3	C	348	0	192	7	1
3	H	285	0	158	4	0
4	A	2	0	0	0	0
4	F	2	0	0	0	0
4	H	1	0	0	0	0
5	F	10	0	0	3	0
6	A	9	0	0	2	0
7	A	159	0	0	11	0
7	B	28	0	0	1	0
7	C	40	0	0	1	0
7	F	166	0	0	11	0
7	G	29	0	0	2	0
7	H	22	0	0	0	0
All	All	7158	0	6221	104	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 8.

All (104) 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:F:129:THR:O	7:F:1001:HOH:O	1.81	0.98
1:F:195:MET:SD	1:F:214:ARG:NH1	2.59	0.76
1:A:193:GLU:OE1	7:A:1001:HOH:O	2.02	0.76
1:A:150:LYS:NZ	3:C:873:DC:OP1	2.18	0.76
1:A:132:ASN:OD1	7:A:1002:HOH:O	2.04	0.75
1:F:157:LYS:HD3	5:F:904:PO4:O1	1.87	0.75
2:G:848:DA:OP2	7:G:901:HOH:O	2.09	0.71
1:F:17:GLU:OE1	7:F:1002:HOH:O	2.07	0.71
1:F:306:ARG:NH2	7:F:1008:HOH:O	2.24	0.70
1:F:127:ARG:NH2	7:F:1006:HOH:O	2.24	0.69
3:C:858:DT:H6	3:C:858:DT:H5"	1.57	0.69
1:A:285:ARG:HB2	1:A:337:THR:HG23	1.77	0.67
1:A:54:ARG:H	1:A:57:MET:HE2	1.59	0.67
1:F:240:ARG:HH21	1:F:277:VAL:HG21	1.60	0.66
1:F:53:VAL:HA	1:F:57:MET:CE	2.26	0.65
1:A:69:LEU:O	7:A:1004:HOH:O	2.13	0.65
1:F:8:ASP:OD2	7:F:1003:HOH:O	2.15	0.64
1:F:53:VAL:HA	1:F:57:MET:HE2	1.80	0.63
1:F:285:ARG:NH1	3:H:866:DT:OP2	2.32	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:VAL:HG23	1:A:173:LEU:HD22	1.82	0.61
2:B:838:DC:N4	7:B:901:HOH:O	2.28	0.61
3:C:858:DT:P	7:C:907:HOH:O	2.59	0.60
2:G:853:DC:OP1	7:G:902:HOH:O	2.17	0.60
1:A:340:ASP:HB2	1:A:341:PRO:HD3	1.84	0.59
1:F:192:LEU:HD23	1:F:195:MET:HE3	1.84	0.58
1:A:247:ARG:NH2	2:B:842:DG:OP1	2.36	0.58
1:F:18:MET:HE1	1:F:45:ASN:HB2	1.86	0.56
1:A:53:VAL:HA	1:A:57:MET:HE1	1.88	0.56
1:F:100:LEU:HD11	1:F:106:TYR:CE2	2.41	0.56
1:F:18:MET:CE	1:F:45:ASN:HD22	2.19	0.55
1:A:53:VAL:HA	1:A:57:MET:CE	2.36	0.55
1:A:24:LEU:HB3	1:A:29:ILE:HG21	1.90	0.53
1:F:18:MET:HE2	1:F:25:ARG:O	2.09	0.53
1:A:157:LYS:HE2	6:A:903:DPO:O5	2.09	0.52
1:A:57:MET:HE3	1:A:62:ALA:HB2	1.92	0.52
1:A:60:GLY:HA3	2:B:838:DC:P	2.49	0.52
1:F:11:CYS:HA	5:F:903:PO4:O2	2.11	0.51
1:F:132:ASN:N	7:F:1001:HOH:O	2.43	0.51
1:A:174:GLN:O	7:A:1005:HOH:O	2.18	0.51
3:C:872:DC:H2'	3:C:873:DC:C6	2.46	0.51
1:F:96:ARG:NH2	7:F:1009:HOH:O	2.33	0.50
5:F:903:PO4:O3	3:H:874:DT:H3'	2.11	0.50
1:A:35:ARG:NH2	7:A:1012:HOH:O	2.35	0.50
1:F:195:MET:SD	1:F:214:ARG:NH2	2.84	0.50
1:F:49:ARG:HD3	7:F:1070:HOH:O	2.12	0.50
1:F:127:ARG:HD2	1:F:138:ALA:O	2.13	0.49
1:A:106:TYR:O	1:A:107:LEU:HD23	2.13	0.49
1:F:267:TYR:OH	1:F:307:LEU:HD13	2.12	0.49
1:A:217:LYS:HG2	2:B:847:DT:OP1	2.12	0.49
1:A:291:LYS:HB3	1:A:331:LEU:HB3	1.94	0.49
1:A:338:LEU:HA	7:A:1003:HOH:O	2.12	0.48
1:F:13:PHE:CD2	3:H:874:DT:H2"	2.49	0.48
1:F:53:VAL:HA	1:F:57:MET:HE1	1.94	0.48
1:A:273:ARG:HD3	7:A:1015:HOH:O	2.13	0.48
1:F:177:PRO:HG2	1:F:180:LYS:HG3	1.94	0.48
1:F:24:LEU:HB3	1:F:29:ILE:HG21	1.94	0.48
1:F:192:LEU:HA	1:F:195:MET:HE2	1.95	0.48
1:A:339:LEU:N	7:A:1003:HOH:O	2.11	0.47
1:F:195:MET:SD	1:F:214:ARG:CZ	3.02	0.47
1:F:18:MET:HE3	1:F:45:ASN:HD22	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:80:LYS:HE3	1:A:84:ASN:HD21	1.79	0.47
1:A:20:ASP:OD2	1:A:75:ARG:NH2	2.46	0.47
1:F:300:GLN:HB3	1:F:323:ARG:NH2	2.31	0.46
1:A:132:ASN:O	1:A:134:LEU:N	2.42	0.46
1:A:278:LYS:HB2	1:A:278:LYS:HE2	1.67	0.46
1:F:207:ASP:OD2	1:F:209:VAL:HB	2.15	0.46
3:C:864:DC:H1'	3:C:865:DC:H5''	1.99	0.45
1:A:100:LEU:HD11	1:A:106:TYR:CE2	2.51	0.45
1:F:187:VAL:O	1:F:191:LYS:HD2	2.16	0.45
1:F:30:ALA:HB2	1:F:69:LEU:HD11	1.98	0.45
1:A:251:GLU:OE1	7:A:1007:HOH:O	2.21	0.45
1:A:269:GLU:OE2	1:A:273:ARG:NE	2.48	0.45
1:A:273:ARG:NH1	7:A:1015:HOH:O	2.45	0.44
1:A:18:MET:CE	1:A:45:ASN:HD22	2.31	0.44
1:A:97:ILE:HG12	1:A:107:LEU:HD22	1.98	0.44
3:C:858:DT:H5''	3:C:858:DT:C6	2.43	0.44
1:A:45:ASN:O	1:A:49:ARG:HG3	2.18	0.44
2:G:848:DA:H2''	2:G:849:DG:C8	2.53	0.43
1:F:85:HIS:CG	1:F:134:LEU:HD11	2.53	0.43
1:F:283:ILE:HD12	1:F:336:VAL:HG22	2.01	0.43
1:F:82:ALA:O	1:F:86:ILE:HG23	2.19	0.43
1:F:57:MET:HE3	1:F:62:ALA:HB2	2.01	0.43
1:A:154:ASP:HA	1:A:157:LYS:HG3	2.01	0.42
1:F:96:ARG:NE	7:F:1009:HOH:O	2.40	0.42
1:F:85:HIS:HB3	1:F:134:LEU:HD21	2.01	0.42
1:F:293:ASP:OD2	1:F:328:GLY:N	2.48	0.42
2:B:842:DG:H2'	2:B:843:DG:O4'	2.19	0.41
1:F:18:MET:HE1	1:F:45:ASN:CB	2.50	0.41
1:F:239:LEU:HA	7:F:1039:HOH:O	2.20	0.41
1:F:32:GLY:HA2	7:F:1068:HOH:O	2.18	0.41
6:A:903:DPO:O7	7:A:1006:HOH:O	2.21	0.41
1:F:54:ARG:H	1:F:57:MET:HE2	1.85	0.41
3:H:872:DC:H2'	3:H:873:DC:C6	2.55	0.41
1:A:5:ILE:HG12	1:A:142:VAL:HG22	2.03	0.41
1:F:173:LEU:O	1:F:200:CYS:HB2	2.19	0.41
1:F:18:MET:HE1	1:F:45:ASN:HD22	1.86	0.41
1:A:240:ARG:HD3	1:A:243:VAL:HG22	2.03	0.41
1:F:18:MET:HG2	1:F:25:ARG:HA	2.03	0.41
1:A:13:PHE:CD2	3:C:874:DT:H2''	2.56	0.41
1:A:192:LEU:HD11	1:A:222:LEU:HD21	2.03	0.40
1:A:289:LYS:HD2	1:A:297:GLN:OE1	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:331:LEU:C	1:A:331:LEU:HD23	2.42	0.40
1:F:178:LEU:HD22	1:F:198:ARG:C	2.42	0.40
1:A:272:ARG:O	1:A:276:LYS:HG2	2.21	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)
1:A:118:SER:OG	3:C:862:DG:OP1[2_455]	2.16	0.04

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	340/352 (97%)	328 (96%)	11 (3%)	1 (0%)	41	47
1	F	340/352 (97%)	328 (96%)	12 (4%)	0	100	100
All	All	680/704 (97%)	656 (96%)	23 (3%)	1 (0%)	51	63

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	129	THR

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	287/297 (97%)	274 (96%)	13 (4%)	27	33
1	F	287/297 (97%)	273 (95%)	14 (5%)	25	29
All	All	574/594 (97%)	547 (95%)	27 (5%)	26	31

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	1	SER
1	F	103	ASP
1	F	154	ASP
1	F	157	LYS
1	F	169	VAL
1	F	178	LEU
1	F	191	LYS
1	F	214	ARG
1	F	222	LEU
1	F	243	VAL
1	F	247	ARG
1	F	260	GLU
1	F	307	LEU
1	F	330	ARG
1	A	1	SER
1	A	35	ARG
1	A	103	ASP
1	A	135	GLN
1	A	154	ASP
1	A	169	VAL
1	A	173	LEU
1	A	191	LYS
1	A	208	LEU
1	A	295	PHE
1	A	336	VAL
1	A	337	THR
1	A	338	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	F	114	HIS
1	F	227	GLN
1	A	84	ASN

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Mol	Chain	Res	Type
1	A	308	ASN

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 5 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
5	PO4	F	904	4	4,4,4	0.84	0	6,6,6	1.74	2 (33%)
5	PO4	F	903	4	4,4,4	1.21	0	6,6,6	1.55	1 (16%)
6	DPO	A	903	4	6,8,8	0.65	0	13,13,13	1.35	1 (7%)

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
6	DPO	A	903	4	-	3/6/6/6	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
6	A	903	DPO	P2-O4-P1	-3.34	121.38	132.83
5	F	904	PO4	O4-P-O1	-3.05	99.74	110.89
5	F	903	PO4	O3-P-O2	-2.99	98.38	107.97
5	F	904	PO4	O4-P-O3	2.43	115.77	107.97

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	903	DPO	P1-O4-P2-O7
6	A	903	DPO	P1-O4-P2-O6
6	A	903	DPO	P1-O4-P2-O5

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	F	904	PO4	1	0
5	F	903	PO4	2	0
6	A	903	DPO	2	0

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	342/352 (97%)	-0.12	4 (1%) 79 86	25, 39, 61, 88	0
1	F	342/352 (97%)	-0.24	0 100 100	20, 33, 55, 74	0
2	B	17/18 (94%)	-0.31	1 (5%) 22 33	32, 44, 57, 96	0
2	G	16/18 (88%)	-0.29	1 (6%) 20 29	23, 36, 79, 90	0
3	C	17/19 (89%)	-0.58	0 100 100	34, 40, 53, 54	0
3	H	14/19 (73%)	-0.48	0 100 100	23, 30, 73, 91	0
All	All	748/778 (96%)	-0.20	6 (0%) 86 91	20, 37, 60, 96	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	0	GLY	4.6
1	A	341	PRO	4.3
2	B	838	DC	3.4
2	G	838	DC	2.7
1	A	76	PHE	2.5
1	A	22	PRO	2.1

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates 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	PO4	F	904	5/5	0.92	0.26	50,57,63,76	0
4	MG	H	901	1/1	0.93	0.47	54,54,54,54	0
4	MG	A	902	1/1	0.97	0.14	49,49,49,49	0
5	PO4	F	903	5/5	0.97	0.21	32,40,50,65	0
6	DPO	A	903	9/9	0.98	0.14	38,44,65,68	0
4	MG	A	901	1/1	0.98	0.09	29,29,29,29	0
4	MG	F	902	1/1	0.98	0.09	37,37,37,37	0
4	MG	F	901	1/1	0.99	0.15	19,19,19,19	0

6.5 Other polymers [i](#)

There are no such residues in this entry.