



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 08:29 pm GMT

PDB ID : 3PX4
Title : Crystal Structure of Bacillus DNA Polymerase I Large Fragment Bound to DNA and ddCTP-dA Mismatch (wobble) in Ajar Conformation
Authors : Wang, W.; Beese, L.S.
Deposited on : 2010-12-09
Resolution : 1.58 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

<http://wwpdb.org/validation/2016/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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

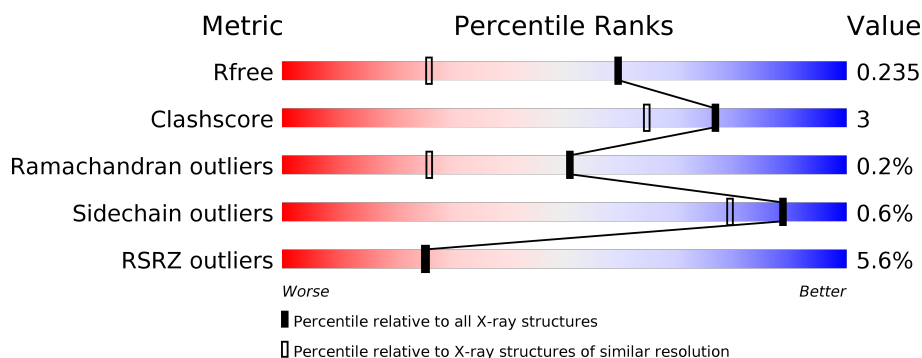
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.58 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}	100719	4211 (1.60-1.56)
Clashscore	112137	4539 (1.60-1.56)
Ramachandran outliers	110173	4423 (1.60-1.56)
Sidechain outliers	110143	4420 (1.60-1.56)
RSRZ outliers	101464	4232 (1.60-1.56)

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. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	592	<div> <div>8%</div> <div>85%</div> <div>8%</div> <div>7%</div> </div>
1	D	592	<div> <div>2%</div> <div>91%</div> <div>7%</div> <div>•</div> </div>
2	B	9	<div> <div>56%</div> <div>33%</div> <div>11%</div> </div>
2	E	9	<div> <div>56%</div> <div>44%</div> </div>
3	C	13	<div> <div>31%</div> <div>38%</div> <div>31%</div> </div>
3	F	13	<div> <div>8%</div> <div>77%</div> <div>15%</div> <div>8%</div> </div>

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 21056 atoms, of which 9648 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called DNA polymerase I.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	549	Total	C	H	N	O	S	0	2	0
			8887	2811	4476	764	821	15			
1	D	579	Total	C	H	N	O	S	0	7	0
			9405	2969	4730	810	879	17			

There are 6 discrepancies between the modelled and reference sequences:

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

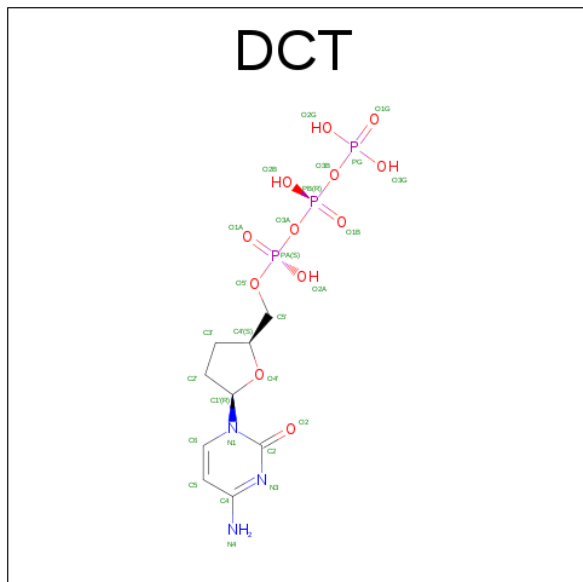
- Molecule 2 is a DNA chain called DNA (5'-D(*CP*CP*TP*GP*AP*CP*TP*CP*(DOC))-3').

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	9	Total	C	H	N	O	P	0	0	0
			275	85	101	29	52	8			
2	E	9	Total	C	H	N	O	P	0	0	0
			275	85	101	29	52	8			

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

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	C	9	Total	C	H	N	O	P	0	0	0
			292	89	101	40	53	9			
3	F	12	Total	C	H	N	O	P	0	0	0
			383	119	133	52	68	11			

- Molecule 4 is 2',3'-DIDEOXYCYTIDINE 5'-TRIPHOSPHATE (three-letter code: DCT) (formula: $C_9H_{16}N_3O_{12}P_3$).





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

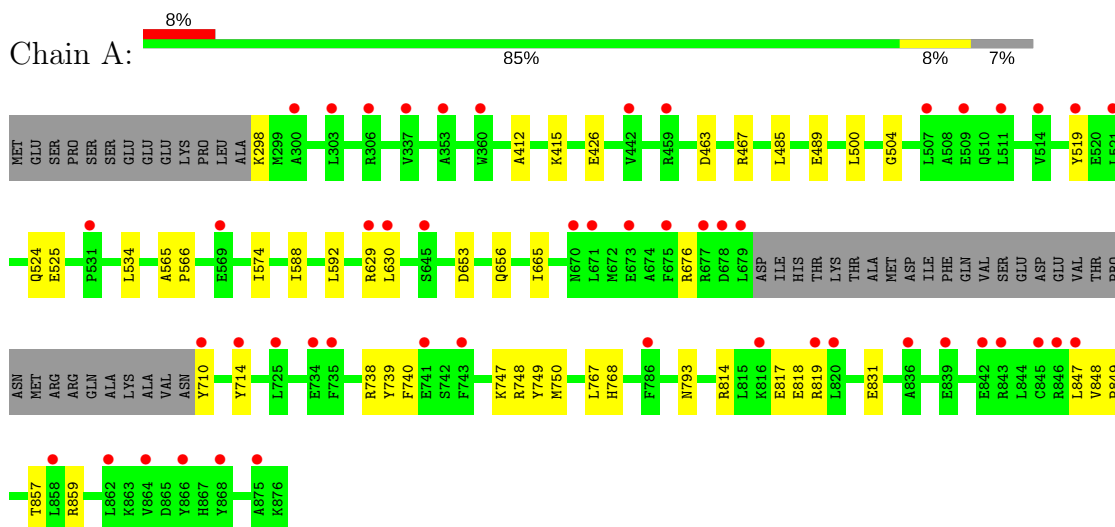
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	503	Total	O	0	0
			503	503		
7	D	775	Total	O	0	0
			775	775		
7	B	31	Total	O	0	0
			31	31		
7	C	50	Total	O	0	0
			50	50		
7	E	41	Total	O	0	0
			41	41		
7	F	68	Total	O	0	0
			68	68		

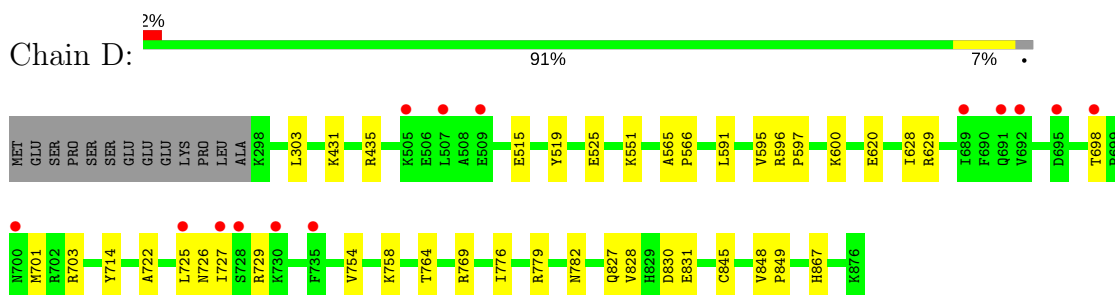
3 Residue-property plots [i](#)

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 I



• Molecule 1: DNA polymerase I



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

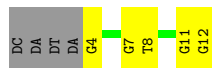


• Molecule 2: DNA (5'-D(*CP*CP*TP*GP*AP*CP*TP*CP*(DOC))-3')

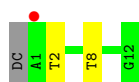
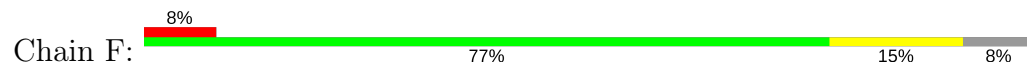




- Molecule 3: DNA (5'-D(*CP*AP*TP*AP*GP*GP*AP*GP*TP*CP*AP*GP*G)-3')



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	93.68Å 109.19Å 150.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	71.10 – 1.58 71.10 – 1.58	Depositor EDS
% Data completeness (in resolution range)	89.6 (71.10-1.58) 89.7 (71.10-1.58)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.90 (at 1.58Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.1_743)	Depositor
R, R_{free}	0.211 , 0.239 0.207 , 0.235	Depositor DCC
R_{free} test set	8577 reflections (4.56%)	DCC
Wilson B-factor (Å ²)	17.7	Xtriage
Anisotropy	0.148	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 48.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	21056	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.76% 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: DOC, MG, DCT, 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	0.35	0/4500	0.50	0/6082
1	D	0.46	0/4782	0.60	0/6464
2	B	0.81	0/173	1.49	3/264 (1.1%)
2	E	0.89	0/173	1.56	2/264 (0.8%)
3	C	0.77	0/215	1.60	6/331 (1.8%)
3	F	1.00	0/282	1.53	4/435 (0.9%)
All	All	0.47	0/10125	0.71	15/13840 (0.1%)

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	23	DT	O4'-C1'-N1	-9.45	101.38	108.00
3	F	8	DT	O4'-C1'-N1	-9.34	101.46	108.00
3	C	8	DT	O4'-C1'-N1	-8.60	101.98	108.00
3	C	7	DG	N1-C6-O6	-7.55	115.37	119.90
3	C	12	DG	O4'-C1'-N9	7.05	112.93	108.00
2	E	27	DT	C4'-C3'-C2'	-6.92	96.87	103.10
3	C	11	DG	O4'-C1'-N9	-6.73	103.29	108.00
2	B	26	DC	O4'-C1'-N1	6.64	112.65	108.00
3	F	2	DT	O4'-C1'-N1	6.62	112.63	108.00
2	B	23	DT	O4'-C1'-N1	-6.39	103.53	108.00
3	C	7	DG	C5-C6-O6	5.37	131.82	128.60
3	F	2	DT	C1'-O4'-C4'	-5.27	104.83	110.10
2	B	27	DT	O4'-C4'-C3'	-5.27	102.39	104.50
3	C	8	DT	N3-C4-O4	5.18	123.01	119.90
3	F	2	DT	O4'-C1'-C2'	-5.11	101.81	105.90

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	4411	4476	4456	33	0
1	D	4675	4730	4686	30	1
2	B	174	101	103	3	0
2	E	174	101	103	1	0
3	C	191	101	101	1	0
3	F	250	133	136	0	0
4	A	27	4	12	0	0
4	D	27	2	12	2	0
5	D	1	0	0	0	0
6	D	10	0	0	0	0
7	A	503	0	0	9	2
7	B	31	0	0	0	0
7	C	50	0	0	1	0
7	D	775	0	0	13	6
7	E	41	0	0	0	0
7	F	68	0	0	0	0
All	All	11408	9648	9609	68	6

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

All (68) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:710:TYR:HB2	7:A:1447:HOH:O	1.89	0.72
1:A:665:ILE:HG22	1:A:749:TYR:CE1	2.28	0.69
1:D:830[B]:ASP:HB3	7:D:969:HOH:O	1.91	0.69
1:D:600:LYS:NZ	7:D:1403:HOH:O	2.30	0.64
1:D:830[A]:ASP:OD1	1:D:831[A]:GLU:HG2	1.99	0.60
1:A:629:ARG:HG3	1:A:630:LEU:HG	1.84	0.59
1:A:738:ARG:NH1	7:A:1108:HOH:O	2.35	0.59
1:D:830[B]:ASP:OD2	4:D:202:DCT:O1A	2.20	0.59
1:A:534:LEU:HD11	1:A:574:ILE:HD13	1.85	0.58
1:A:740:PHE:HB3	1:A:747:LYS:HD2	1.85	0.58
1:D:725:LEU:O	1:D:727:ILE:HG23	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:748:ARG:HD2	7:A:1245:HOH:O	2.04	0.57
1:A:519:TYR:CD2	1:A:525:GLU:HG2	2.41	0.56
1:D:828:VAL:HB	1:D:831[B]:GLU:CG	2.35	0.56
1:A:848:VAL:HB	1:A:849:PRO:HD3	1.87	0.56
1:D:776:ILE:HG22	7:D:1129:HOH:O	2.05	0.55
1:D:629:ARG:NH1	7:D:1271:HOH:O	2.39	0.55
1:D:565:ALA:N	1:D:566:PRO:CD	2.70	0.54
1:A:714:TYR:HB2	7:A:1018:HOH:O	2.08	0.53
1:D:515:GLU:HG2	1:D:519:TYR:CE2	2.43	0.53
1:D:629:ARG:NH2	7:D:1216:HOH:O	2.34	0.53
1:A:504:GLY:HA2	1:A:588:ILE:HD12	1.89	0.52
1:D:551:LYS:NZ	7:D:1046:HOH:O	2.42	0.52
2:B:22:DC:H5'	2:B:22:DC:C6	2.45	0.51
1:D:435:ARG:NH2	7:D:1463:HOH:O	2.35	0.51
1:A:629:ARG:HD2	7:A:1131:HOH:O	2.10	0.50
1:D:726:ASN:CG	7:D:1433:HOH:O	2.50	0.50
1:D:828:VAL:HB	1:D:831[A]:GLU:HG3	1.94	0.50
1:D:776:ILE:CG2	7:D:1129:HOH:O	2.57	0.49
1:A:412:ALA:HA	1:A:415:LYS:HD3	1.95	0.49
1:D:714:TYR:HB3	7:D:908:HOH:O	2.13	0.49
1:D:845:CYS:SG	7:D:883:HOH:O	2.48	0.48
1:A:749:TYR:HD2	1:A:750:MET:HE3	1.78	0.48
1:A:767:LEU:O	1:A:768:HIS:HB2	2.13	0.48
1:D:620:GLU:HB2	7:D:1209:HOH:O	2.13	0.48
2:B:22:DC:H5'	2:B:22:DC:H6	1.78	0.48
1:A:298:LYS:N	7:A:1036:HOH:O	2.47	0.47
1:D:725:LEU:HB2	1:D:727:ILE:HG12	1.96	0.47
1:A:676:ARG:NH2	1:A:859:ARG:O	2.47	0.47
1:D:848:VAL:HB	1:D:849:PRO:HD3	1.97	0.47
1:A:485:LEU:O	1:A:489:GLU:HG3	2.14	0.46
1:A:426:GLU:HG3	7:A:1258:HOH:O	2.14	0.46
1:A:565:ALA:N	1:A:566:PRO:CD	2.78	0.46
1:A:817:GLU:HG3	1:A:818:GLU:N	2.30	0.45
1:D:519:TYR:CD2	1:D:525:GLU:HG2	2.50	0.45
1:A:814:ARG:HB3	1:A:847:LEU:HD11	1.97	0.45
1:D:867:HIS:HD2	7:D:1205:HOH:O	2.00	0.45
1:A:463:ASP:O	1:A:467:ARG:HG3	2.17	0.45
1:D:698:THR:OG1	1:D:701:MET:HG3	2.18	0.44
1:A:500:LEU:HD23	1:A:592:LEU:HG	1.99	0.44
1:A:665:ILE:CG2	1:A:749:TYR:CE1	2.98	0.43
1:A:814:ARG:O	1:A:818:GLU:HG2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:793:ASN:HB2	7:A:1046:HOH:O	2.16	0.43
4:D:202:DCT:O5'	4:D:202:DCT:H6	2.18	0.43
1:D:722:ALA:HB2	1:D:729:ARG:HA	2.00	0.43
1:A:817:GLU:HG3	1:A:818:GLU:H	1.83	0.43
1:D:754:VAL:CG1	1:D:758:LYS:HE2	2.49	0.43
3:C:4:DG:H5''	7:C:215:HOH:O	2.18	0.43
1:A:656:GLN:NE2	7:A:1417:HOH:O	2.36	0.42
1:D:764:THR:HA	1:D:769:ARG:O	2.19	0.42
1:D:591:LEU:O	1:D:595:VAL:HG23	2.20	0.41
1:D:596:ARG:HA	1:D:597:PRO:HD3	1.96	0.41
1:D:754:VAL:HG12	1:D:758:LYS:HE2	2.02	0.41
2:E:28:DC:H2'	2:E:29:DOC:H6	2.02	0.41
1:A:524:GLN:HG2	1:A:525:GLU:N	2.36	0.41
1:A:739:TYR:C	1:A:739:TYR:CD2	2.94	0.41
1:A:653:ASP:CB	1:A:831:GLU:HG2	2.51	0.41
1:A:629:ARG:HH12	2:B:27:DT:P	2.45	0.40

All (6) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:D:1210:HOH:O	7:D:1238:HOH:O[4_445]	1.69	0.51
7:A:1211:HOH:O	7:D:1208:HOH:O[2_745]	1.82	0.38
1:D:779:ARG:HH22	7:D:1406:HOH:O[4_545]	1.39	0.21
7:D:1027:HOH:O	7:D:1028:HOH:O[4_545]	2.03	0.17
7:A:990:HOH:O	7:D:942:HOH:O[4_545]	2.14	0.06
7:D:1397:HOH:O	7:D:1407:HOH:O[4_445]	2.17	0.03

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	547/592 (92%)	531 (97%)	15 (3%)	1 (0%)	51	26
1	D	584/592 (99%)	571 (98%)	12 (2%)	1 (0%)	51	26
All	All	1131/1184 (96%)	1102 (97%)	27 (2%)	2 (0%)	51	26

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	819	ARG
1	D	628	ILE

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	469/507 (92%)	468 (100%)	1 (0%)	94	89
1	D	501/507 (99%)	496 (99%)	5 (1%)	80	63
All	All	970/1014 (96%)	964 (99%)	6 (1%)	89	79

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

Mol	Chain	Res	Type
1	A	857	THR
1	D	303	LEU
1	D	431	LYS
1	D	703	ARG
1	D	782	ASN
1	D	827	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	543	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	DOC	B	29	3,2	13,19,20	1.03	1 (7%)	12,26,29	0.74	0
2	DOC	E	29	3,2	13,19,20	1.09	1 (7%)	12,26,29	0.63	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	DOC	B	29	3,2	-	0/3/18/19	0/2/2/2
2	DOC	E	29	3,2	-	0/3/18/19	0/2/2/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	29	DOC	O5'-C5'	-2.82	1.40	1.44
2	B	29	DOC	C2-N3	-2.22	1.33	1.38

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	29	DOC	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	DCT	A	203	-	22,28,28	0.78	0	20,43,43	1.58	5 (25%)
6	SO4	D	2	-	4,4,4	0.23	0	6,6,6	0.10	0
4	DCT	D	202	5	22,28,28	0.83	1 (4%)	20,43,43	1.23	1 (5%)
6	SO4	D	877	-	4,4,4	0.17	0	6,6,6	0.31	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	DCT	A	203	-	-	0/18/31/31	0/2/2/2
6	SO4	D	2	-	-	0/0/0/0	0/0/0/0
4	DCT	D	202	5	-	0/18/31/31	0/2/2/2
6	SO4	D	877	-	-	0/0/0/0	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	202	DCT	PG-O3B	2.01	1.63	1.60

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	202	DCT	C2'-C1'-N1	-3.68	105.46	112.47
4	A	203	DCT	C2'-C1'-N1	-3.02	106.71	112.47
4	A	203	DCT	C5-C4-N3	-2.01	119.28	121.68
4	A	203	DCT	N4-C4-N3	2.05	120.10	116.64
4	A	203	DCT	C3'-C2'-C1'	2.89	105.90	102.69
4	A	203	DCT	O4'-C4'-C5'	3.60	115.43	109.59

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	202	DCT	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	549/592 (92%)	0.52	50 (9%) 10 10	14, 31, 50, 64	0
1	D	579/592 (97%)	0.18	14 (2%) 59 61	7, 18, 38, 53	0
2	B	8/9 (88%)	-0.23	0 100 100	19, 24, 42, 47	0
2	E	8/9 (88%)	-0.18	0 100 100	15, 22, 38, 48	0
3	C	9/13 (69%)	-0.42	0 100 100	17, 20, 27, 33	0
3	F	12/13 (92%)	0.27	1 (8%) 12 12	12, 20, 44, 57	0
All	All	1165/1228 (94%)	0.34	65 (5%) 25 25	7, 24, 47, 64	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	819	ARG	5.8
1	A	866	TYR	5.7
1	A	679	LEU	5.4
1	A	671	LEU	4.8
1	A	868	TYR	4.6
1	A	847	LEU	4.5
1	A	842	GLU	4.2
1	A	843	ARG	4.1
1	A	630	LEU	4.0
1	A	846	ARG	4.0
1	A	303	LEU	3.8
1	D	509	GLU	3.8
1	A	710	TYR	3.7
1	A	862	LEU	3.7
1	A	743	PHE	3.6
1	A	677	ARG	3.4
1	A	521	LEU	3.0
1	A	864	VAL	2.9
1	D	692	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	507	LEU	2.8
1	A	360	TRP	2.8
1	A	675	PHE	2.8
1	A	300	ALA	2.8
1	A	816	LYS	2.8
1	A	645	SER	2.8
1	A	670	ASN	2.8
1	A	714	TYR	2.8
1	A	820	LEU	2.7
1	A	629	ARG	2.7
1	D	725	LEU	2.7
1	D	698	THR	2.7
1	D	735	PHE	2.7
1	A	507	LEU	2.7
1	A	725	LEU	2.6
1	D	691	GLN	2.6
1	A	306	ARG	2.6
1	A	678	ASP	2.5
1	A	531	PRO	2.5
1	A	858	LEU	2.4
1	D	505	LYS	2.4
1	A	786	PHE	2.4
1	D	695	ASP	2.4
1	A	459	ARG	2.4
1	D	700	ASN	2.3
1	A	442	VAL	2.3
1	A	569	GLU	2.3
1	A	337	VAL	2.3
1	A	845	CYS	2.3
1	A	353	ALA	2.3
1	D	689	ILE	2.3
1	D	727	ILE	2.3
1	A	519	TYR	2.3
3	F	1	DA	2.3
1	D	728	SER	2.2
1	A	875	ALA	2.2
1	A	735	PHE	2.2
1	A	836	ALA	2.2
1	D	730	LYS	2.2
1	A	734	GLU	2.2
1	A	511	LEU	2.2
1	A	839	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	509	GLU	2.1
1	A	741	GLU	2.1
1	A	673	GLU	2.1
1	A	514	VAL	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	DOC	B	29	18/19	0.96	0.08	-	15,24,32,35	0
2	DOC	E	29	18/19	0.98	0.11	-	10,14,19,22	0

6.3 Carbohydrates [i](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
6	SO4	D	2	5/5	0.94	0.09	-0.05	27,29,32,38	0
4	DCT	A	203	27/27	0.90	0.09	-0.13	19,36,62,65	0
4	DCT	D	202	27/27	0.97	0.08	-0.60	14,20,29,31	0
5	MG	D	1	1/1	0.98	0.06	-2.09	29,29,29,29	0
6	SO4	D	877	5/5	0.99	0.09	-	26,26,37,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.