



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:41 PM GMT

PDB ID : 4FS2
Title : Base pairing mechanism of N2,3-ethenoguanine with dCTP by human polymerase ι
Authors : Zhao, L.
Deposited on : 2012-06-26
Resolution : 2.05 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

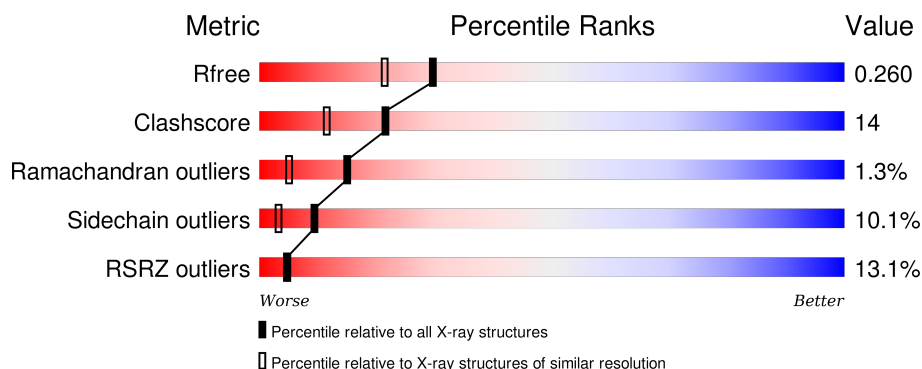
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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}	91344	1192 (2.04-2.04)
Clashscore	102246	1269 (2.04-2.04)
Ramachandran outliers	100387	1258 (2.04-2.04)
Sidechain outliers	100360	1258 (2.04-2.04)
RSRZ outliers	91569	1194 (2.04-2.04)

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	B	18	
1	C	18	
2	A	420	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3422 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 DNA chain called DNA (5'-D(*TP*CP*TP*(EFG)P*GP*GP*GP*TP*CP*CP*TP*AP*GP*GP*AP*CP*CP*(DOC))-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	7	Total	C	N	O	P	0	0	0
			142	67	29	39	7			
1	C	8	Total	C	F	N	O	P	0	0
			169	80	1	30	50	8		

- Molecule 2 is a protein called DNA polymerase iota.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	383	Total	C	N	O	S	0	0	0
			2900	1828	511	540	21			

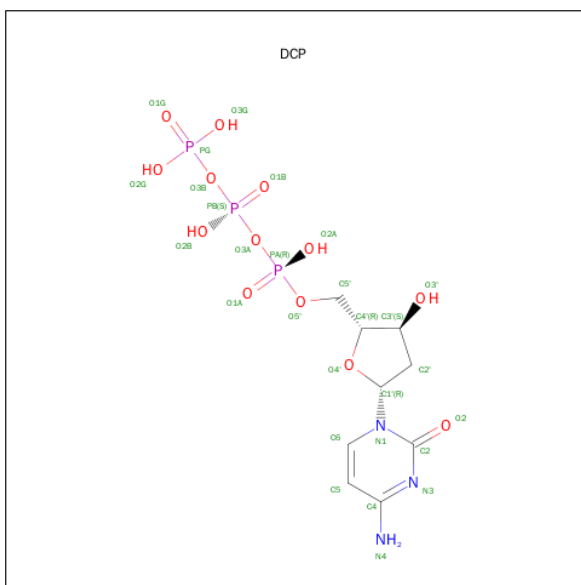
- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total	Mg	0	0
			3	3		

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

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

- Molecule 5 is 2'-DEOXYCYTIDINE-5'-TRIPHOSPHATE (three-letter code: DCP) (formula: C₉H₁₆N₃O₁₃P₃).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	8	Total O 8 8	0	0
6	C	22	Total O 22 22	0	0
6	A	149	Total O 149 149	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	97.30 Å 97.30 Å 202.91 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	84.26 – 2.05 28.81 – 2.05	Depositor EDS
% Data completeness (in resolution range)	99.7 (84.26-2.05) 99.8 (28.81-2.05)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.31 (at 2.04 Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.214 , 0.261 0.218 , 0.260	Depositor DCC
R_{free} test set	1822 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	42.8	Xtriage
Anisotropy	0.145	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 59.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 36362 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3422	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.44% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: NA, MG, EFG, DCP, DOC

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	B	0.60	0/139	0.99	0/212
1	C	1.36	2/160 (1.2%)	0.90	1/245 (0.4%)
2	A	0.89	0/2940	0.94	4/3982 (0.1%)
All	All	0.91	2/3239 (0.1%)	0.94	5/4439 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	842	DG	O3'-P	-6.66	1.53	1.61
1	C	843	DG	O3'-P	-5.99	1.53	1.61

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	845	DC	P-O3'-C3'	6.97	128.06	119.70
2	A	135	MET	CG-SD-CE	-5.50	91.39	100.20
2	A	368	VAL	CB-CA-C	-5.50	100.94	111.40
2	A	344	LEU	CA-CB-CG	5.18	127.23	115.30
2	A	413	LEU	CA-CB-CG	5.05	126.91	115.30

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	B	142	0	78	3	2
1	C	169	0	90	5	2
2	A	2900	0	2897	83	0
3	A	3	0	0	0	0
4	A	1	0	0	0	0
5	A	28	0	12	0	0
6	A	149	0	0	5	0
6	B	8	0	0	0	0
6	C	22	0	0	0	0
All	All	3422	0	3077	89	2

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 14.

All (89) 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)
2:A:310:LYS:HE2	2:A:349:TYR:CB	1.74	1.18
2:A:237:LYS:HE2	6:A:746:HOH:O	1.42	1.16
2:A:35:LEU:HD21	2:A:187:MET:HE1	1.31	1.13
2:A:137:GLU:HG2	2:A:172:ARG:NH2	1.75	1.01
2:A:310:LYS:CE	2:A:349:TYR:CB	2.43	0.96
1:B:11:DC:OP1	2:A:246:THR:HG22	1.67	0.95
2:A:275:ILE:HD11	2:A:279:GLN:HE21	1.31	0.93
2:A:275:ILE:HD11	2:A:279:GLN:NE2	1.84	0.92
2:A:35:LEU:HD21	2:A:187:MET:CE	2.02	0.90
2:A:164:ASN:H	2:A:170:HIS:HD2	1.19	0.89
2:A:137:GLU:CG	2:A:172:ARG:HH21	1.89	0.86
2:A:137:GLU:HG2	2:A:172:ARG:HH21	1.36	0.85
2:A:103:ARG:CZ	2:A:331:ARG:HD3	2.08	0.84
1:C:844:DT:H2''	1:C:845:DC:H5'	1.59	0.82
2:A:275:ILE:CD1	2:A:279:GLN:HE21	1.97	0.77
1:C:841:DG:OP2	2:A:307:SER:HB2	1.86	0.76
2:A:331:ARG:HD2	6:A:747:HOH:O	1.85	0.76
2:A:144:GLN:CD	2:A:144:GLN:H	1.91	0.71
2:A:266:PRO:O	2:A:270:GLU:HG3	1.93	0.68
2:A:164:ASN:H	2:A:170:HIS:CD2	2.09	0.67
2:A:314:GLU:O	2:A:316:GLU:N	2.31	0.64
2:A:112:LEU:HD23	2:A:112:LEU:C	2.18	0.64
2:A:318:LYS:HE3	2:A:384:VAL:HG11	1.81	0.63
1:C:840:EFG:H10	1:C:840:EFG:O4'	1.99	0.61
2:A:345:ILE:HB	2:A:407:SER:HB3	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:243:GLY:H	2:A:246:THR:CG2	2.14	0.60
2:A:368:VAL:O	2:A:368:VAL:HG12	2.00	0.60
2:A:343:ARG:HD3	2:A:345:ILE:HD11	1.84	0.58
2:A:369:ILE:HG22	2:A:369:ILE:O	2.02	0.58
2:A:304:GLU:HG2	2:A:331:ARG:HH21	1.67	0.58
2:A:144:GLN:N	2:A:144:GLN:CD	2.57	0.57
2:A:100:THR:O	2:A:104:GLU:HG3	2.05	0.57
2:A:103:ARG:NH2	2:A:331:ARG:HD3	2.20	0.56
2:A:164:ASN:N	2:A:170:HIS:HD2	1.98	0.55
1:C:844:DT:H2''	1:C:845:DC:C5'	2.35	0.55
2:A:316:GLU:HG2	2:A:317:ALA:N	2.23	0.54
2:A:303:SER:O	2:A:331:ARG:NH2	2.40	0.54
2:A:348:ARG:NH1	2:A:394:MET:O	2.36	0.54
2:A:368:VAL:O	2:A:370:GLN:N	2.40	0.54
2:A:379:VAL:O	2:A:382:PRO:HD2	2.09	0.53
2:A:242:ILE:HA	2:A:246:THR:HG21	1.90	0.52
2:A:144:GLN:OE1	2:A:147:GLU:CB	2.58	0.52
2:A:35:LEU:CD2	2:A:187:MET:CE	2.82	0.51
2:A:243:GLY:N	2:A:246:THR:CG2	2.73	0.51
2:A:178:GLN:HG2	6:A:717:HOH:O	2.11	0.51
2:A:320:LYS:HE3	2:A:320:LYS:HA	1.91	0.51
2:A:320:LYS:HE3	2:A:320:LYS:CA	2.40	0.51
2:A:325:LEU:CD1	2:A:329:LEU:HD22	2.42	0.50
2:A:202:ASN:ND2	2:A:205:LEU:H	2.08	0.50
2:A:309:LYS:O	2:A:311:CYS:N	2.44	0.50
2:A:304:GLU:HG3	2:A:328:LEU:HG	1.92	0.50
2:A:412:ASN:HD21	2:A:414:LYS:HE3	1.76	0.50
2:A:316:GLU:O	2:A:319:ASN:N	2.45	0.50
2:A:371:LYS:HD2	2:A:371:LYS:N	2.27	0.50
1:B:9:DG:H2''	1:B:10:DA:C8	2.47	0.49
2:A:103:ARG:NH1	2:A:331:ARG:HD3	2.27	0.48
2:A:325:LEU:HD12	2:A:380:MET:SD	2.52	0.48
2:A:88:CYS:O	2:A:91:LEU:HB2	2.12	0.48
2:A:412:ASN:ND2	2:A:414:LYS:HE3	2.29	0.47
2:A:331:ARG:CD	6:A:747:HOH:O	2.52	0.47
2:A:265:SER:OG	2:A:268:ILE:HG13	2.15	0.47
1:B:10:DA:H2'	1:B:11:DC:C6	2.49	0.47
2:A:138:LYS:O	2:A:142:GLN:HG2	2.14	0.47
2:A:343:ARG:HG2	2:A:344:LEU:N	2.30	0.46
1:C:844:DT:H2'	1:C:845:DC:C6	2.51	0.45
2:A:202:ASN:C	2:A:202:ASN:HD22	2.20	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:304:GLU:HG2	2:A:331:ARG:NH2	2.32	0.45
2:A:380:MET:O	2:A:384:VAL:HG23	2.17	0.45
2:A:243:GLY:H	2:A:246:THR:HG22	1.82	0.45
2:A:325:LEU:HD13	2:A:325:LEU:O	2.16	0.45
2:A:62:LEU:HA	2:A:62:LEU:HD22	1.73	0.44
2:A:313:SER:OG	2:A:401:PHE:CB	2.65	0.44
2:A:310:LYS:HZ2	2:A:402:HIS:HA	1.82	0.44
2:A:379:VAL:C	2:A:382:PRO:HD2	2.39	0.43
2:A:61:TYR:CG	2:A:61:TYR:O	2.70	0.43
2:A:105:MET:CG	2:A:193:LEU:HD11	2.49	0.43
2:A:325:LEU:HD13	2:A:325:LEU:C	2.40	0.42
2:A:253:LEU:HD22	2:A:264:PHE:CZ	2.54	0.42
2:A:253:LEU:HD22	2:A:264:PHE:HZ	1.85	0.41
2:A:140:LEU:HD23	2:A:140:LEU:HA	1.89	0.41
2:A:112:LEU:CD2	2:A:112:LEU:C	2.88	0.41
2:A:365:PRO:HB2	2:A:368:VAL:HG23	2.02	0.41
2:A:105:MET:HE1	6:A:741:HOH:O	2.20	0.41
2:A:310:LYS:HA	2:A:310:LYS:HD3	1.80	0.41
2:A:368:VAL:CG1	2:A:368:VAL:O	2.67	0.41
2:A:61:TYR:CD1	2:A:62:LEU:HD23	2.55	0.41
2:A:329:LEU:O	2:A:333:CYS:HB2	2.20	0.40
2:A:55:LEU:HG	2:A:56:GLY:N	2.37	0.40
2:A:58:GLN:HA	2:A:62:LEU:O	2.22	0.40

All (2) 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:B:7:DA:P	1:C:847:DT:O3'[10_445]	1.35	0.85
1:B:7:DA:O5'	1:C:847:DT:O3'[10_445]	2.15	0.05

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
2	A	379/420 (90%)	358 (94%)	16 (4%)	5 (1%)	15 5

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	310	LYS
2	A	315	VAL
2	A	317	ALA
2	A	73	LEU
2	A	369	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
2	A	318/376 (85%)	286 (90%)	32 (10%)	9 3

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	27	SER
2	A	61	TYR
2	A	62	LEU
2	A	80	ASN
2	A	91	LEU
2	A	126	ASP
2	A	144	GLN
2	A	149	SER
2	A	160	ASN
2	A	170	HIS
2	A	202	ASN
2	A	265	SER
2	A	290	ASN
2	A	294	ILE
2	A	295	LEU

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Mol	Chain	Res	Type
2	A	305	GLU
2	A	307	SER
2	A	310	LYS
2	A	311	CYS
2	A	312	SER
2	A	313	SER
2	A	316	GLU
2	A	318	LYS
2	A	320	LYS
2	A	329	LEU
2	A	333	CYS
2	A	343	ARG
2	A	344	LEU
2	A	357	ARG
2	A	370	GLN
2	A	398	LYS
2	A	413	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	170	HIS
2	A	202	ASN
2	A	262	GLN
2	A	279	GLN
2	A	330	ASN
2	A	340	HIS
2	A	412	ASN

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	$\# Z > 2$
1	DOC	B	13	1	11,19,20	0.91	1 (9%)	14,26,29	1.47	4 (28%)
1	EFG	C	840	-	16,28,29	2.21	5 (31%)	17,42,45	2.38	7 (41%)

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
1	DOC	B	13	1	-	0/3/18/19	0/2/2/2
1	EFG	C	840	-	-	0/3/25/26	0/4/4/4

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	840	EFG	C6-C5	-4.39	1.32	1.41
1	C	840	EFG	C2-N2	-3.77	1.30	1.35
1	C	840	EFG	O3'-C3'	-3.23	1.35	1.43
1	C	840	EFG	C5-C4	-3.11	1.33	1.40
1	B	13	DOC	C6-N1	-2.24	1.32	1.35
1	C	840	EFG	C2'-C3'	-2.12	1.49	1.52

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	840	EFG	C5-C6-N1	-4.97	116.79	123.59
1	C	840	EFG	C4'-O4'-C1'	-4.92	104.31	109.72
1	C	840	EFG	O4'-C1'-C2'	-2.86	102.60	105.84
1	C	840	EFG	C4-C5-N7	-2.84	106.87	109.48
1	C	840	EFG	F-C2'-C3'	-2.57	102.82	109.07
1	C	840	EFG	O3'-C3'-C4'	-2.51	103.54	111.05
1	B	13	DOC	C2-N3-C4	2.13	118.62	115.61
1	C	840	EFG	O4'-C4'-C5'	2.18	117.12	109.32
1	B	13	DOC	O4'-C1'-N1	2.28	111.66	107.72
1	B	13	DOC	C3'-C2'-C1'	2.35	105.34	102.71
1	B	13	DOC	O4'-C4'-C3'	2.68	109.18	104.69

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
1	C	840	EFG	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, 4 are monoatomic - leaving 1 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	DCP	A	503	3	21,29,29	1.06	1 (4%)	33,45,45	1.64	5 (15%)

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
5	DCP	A	503	3	-	0/18/34/34	0/2/2/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	503	DCP	C5-C4	-2.30	1.35	1.40

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	503	DCP	C2'-C1'-N1	-2.14	108.97	114.16
5	A	503	DCP	C5-C4-N3	-2.03	119.23	121.80
5	A	503	DCP	O2A-PA-O1A	2.39	125.50	112.53
5	A	503	DCP	O3G-PG-O2G	3.00	118.79	107.38
5	A	503	DCP	C2-N3-C4	5.94	124.00	115.61

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	B	6/18 (33%)	-0.34	0	100 100	42, 56, 59, 63	0
1	C	7/18 (38%)	-0.49	0	100 100	36, 39, 48, 49	0
2	A	383/420 (91%)	0.73	52 (13%)	4 4	28, 50, 94, 176	1 (0%)
All	All	396/456 (86%)	0.69	52 (13%)	5 5	28, 50, 94, 176	1 (0%)

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	353	LYS	15.5
2	A	355	TYR	13.7
2	A	352	GLU	11.7
2	A	354	HIS	10.9
2	A	311	CYS	9.4
2	A	336	GLY	5.2
2	A	400	PRO	4.9
2	A	84	ALA	4.8
2	A	351	SER	4.5
2	A	82	ARG	4.3
2	A	142	GLN	4.2
2	A	398	LYS	3.6
2	A	33	VAL	3.6
2	A	252	ALA	3.6
2	A	333	CYS	3.5
2	A	78	LEU	3.4
2	A	397	VAL	3.3
2	A	80	ASN	3.3
2	A	350	SER	3.3
2	A	26	SER	3.2
2	A	61	TYR	3.1
2	A	62	LEU	3.0
2	A	109	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
2	A	268	ILE	2.9
2	A	73	LEU	2.9
2	A	130	VAL	2.9
2	A	79	MET	2.9
2	A	31	VAL	2.8
2	A	81	VAL	2.8
2	A	378	ASP	2.8
2	A	75	VAL	2.8
2	A	143	LEU	2.7
2	A	348	ARG	2.7
2	A	356	GLY	2.7
2	A	399	MET	2.6
2	A	113	LEU	2.6
2	A	371	LYS	2.5
2	A	387	LEU	2.5
2	A	86	GLU	2.5
2	A	83	ASP	2.5
2	A	91	LEU	2.4
2	A	335	ASP	2.4
2	A	367	HIS	2.4
2	A	145	SER	2.4
2	A	269	LEU	2.3
2	A	129	PHE	2.3
2	A	144	GLN	2.1
2	A	106	SER	2.1
2	A	264	PHE	2.1
2	A	123	LEU	2.1
2	A	51	LYS	2.1
2	A	342	VAL	2.0

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

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
1	EFG	C	840	25/26	0.97	0.09	-1.06	40,48,67,71	0
1	DOC	B	13	18/19	0.99	0.15	-	40,44,48,48	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(\AA^2)	Q<0.9
3	MG	A	501	1/1	0.97	0.24	1.98	15,15,15,15	1
3	MG	A	504	1/1	0.96	0.25	1.72	41,41,41,41	0
5	DCP	A	503	28/28	0.98	0.11	-0.80	29,38,42,44	0
4	NA	A	502	1/1	0.96	0.06	-1.51	38,38,38,38	0
3	MG	A	505	1/1	0.86	0.87	-	30,30,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.