



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:27 AM GMT

PDB ID : 2QRN
Title : Human Deoxycytidine kinase dCMP, UDP, Mg ion product complex
Authors : Ealick, S.E.; Soriano, E.V.
Deposited on : 2007-07-28
Resolution : 3.40 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

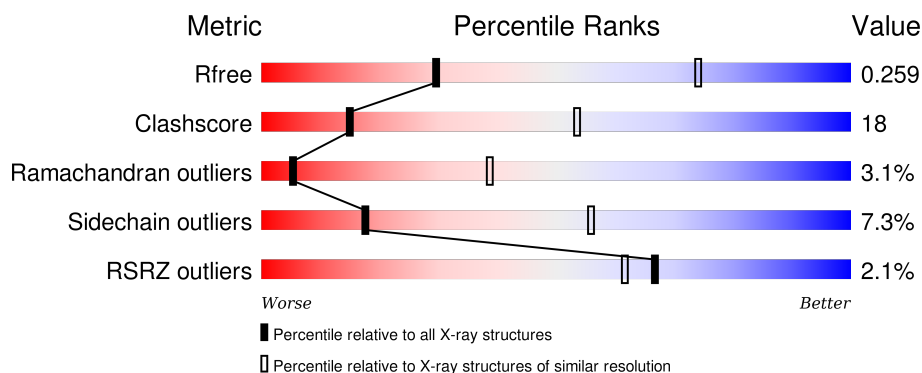
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 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	1476 (3.50-3.30)
Clashscore	102246	1611 (3.50-3.30)
Ramachandran outliers	100387	1571 (3.50-3.30)
Sidechain outliers	100360	1571 (3.50-3.30)
RSRZ outliers	91569	1485 (3.50-3.30)

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	280	<div> <div>3%</div> <div>52%</div> <div>24%</div> <div>5%</div> <div>19%</div> </div>
1	B	280	<div> <div>%</div> <div>53%</div> <div>24%</div> <div>.</div> <div>19%</div> </div>
1	C	280	<div> <div>3%</div> <div>52%</div> <div>23%</div> <div>6%</div> <div>19%</div> </div>
1	D	280	<div> <div>%</div> <div>53%</div> <div>24%</div> <div>.</div> <div>19%</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
4	DCM	C	403	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7808 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 Deoxycytidine kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	228	Total	C	N	O	S	0	0	0
			1906	1223	317	357	9			
1	B	228	Total	C	N	O	S	0	0	0
			1906	1223	317	357	9			
1	C	228	Total	C	N	O	S	0	0	0
			1906	1223	317	357	9			
1	D	228	Total	C	N	O	S	0	0	0
			1906	1223	317	357	9			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP P27707
A	-18	GLY	-	expression tag	UNP P27707
A	-17	SER	-	expression tag	UNP P27707
A	-16	SER	-	expression tag	UNP P27707
A	-15	HIS	-	expression tag	UNP P27707
A	-14	HIS	-	expression tag	UNP P27707
A	-13	HIS	-	expression tag	UNP P27707
A	-12	HIS	-	expression tag	UNP P27707
A	-11	HIS	-	expression tag	UNP P27707
A	-10	HIS	-	expression tag	UNP P27707
A	-9	SER	-	expression tag	UNP P27707
A	-8	SER	-	expression tag	UNP P27707
A	-7	GLY	-	expression tag	UNP P27707
A	-6	LEU	-	expression tag	UNP P27707
A	-5	VAL	-	expression tag	UNP P27707
A	-4	PRO	-	expression tag	UNP P27707
A	-3	ARG	-	expression tag	UNP P27707
A	-2	GLY	-	expression tag	UNP P27707
A	-1	SER	-	expression tag	UNP P27707
A	0	HIS	-	expression tag	UNP P27707
B	-19	MET	-	expression tag	UNP P27707

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	GLY	-	expression tag	UNP P27707
B	-17	SER	-	expression tag	UNP P27707
B	-16	SER	-	expression tag	UNP P27707
B	-15	HIS	-	expression tag	UNP P27707
B	-14	HIS	-	expression tag	UNP P27707
B	-13	HIS	-	expression tag	UNP P27707
B	-12	HIS	-	expression tag	UNP P27707
B	-11	HIS	-	expression tag	UNP P27707
B	-10	HIS	-	expression tag	UNP P27707
B	-9	SER	-	expression tag	UNP P27707
B	-8	SER	-	expression tag	UNP P27707
B	-7	GLY	-	expression tag	UNP P27707
B	-6	LEU	-	expression tag	UNP P27707
B	-5	VAL	-	expression tag	UNP P27707
B	-4	PRO	-	expression tag	UNP P27707
B	-3	ARG	-	expression tag	UNP P27707
B	-2	GLY	-	expression tag	UNP P27707
B	-1	SER	-	expression tag	UNP P27707
B	0	HIS	-	expression tag	UNP P27707
C	-19	MET	-	expression tag	UNP P27707
C	-18	GLY	-	expression tag	UNP P27707
C	-17	SER	-	expression tag	UNP P27707
C	-16	SER	-	expression tag	UNP P27707
C	-15	HIS	-	expression tag	UNP P27707
C	-14	HIS	-	expression tag	UNP P27707
C	-13	HIS	-	expression tag	UNP P27707
C	-12	HIS	-	expression tag	UNP P27707
C	-11	HIS	-	expression tag	UNP P27707
C	-10	HIS	-	expression tag	UNP P27707
C	-9	SER	-	expression tag	UNP P27707
C	-8	SER	-	expression tag	UNP P27707
C	-7	GLY	-	expression tag	UNP P27707
C	-6	LEU	-	expression tag	UNP P27707
C	-5	VAL	-	expression tag	UNP P27707
C	-4	PRO	-	expression tag	UNP P27707
C	-3	ARG	-	expression tag	UNP P27707
C	-2	GLY	-	expression tag	UNP P27707
C	-1	SER	-	expression tag	UNP P27707
C	0	HIS	-	expression tag	UNP P27707
D	-19	MET	-	expression tag	UNP P27707
D	-18	GLY	-	expression tag	UNP P27707
D	-17	SER	-	expression tag	UNP P27707

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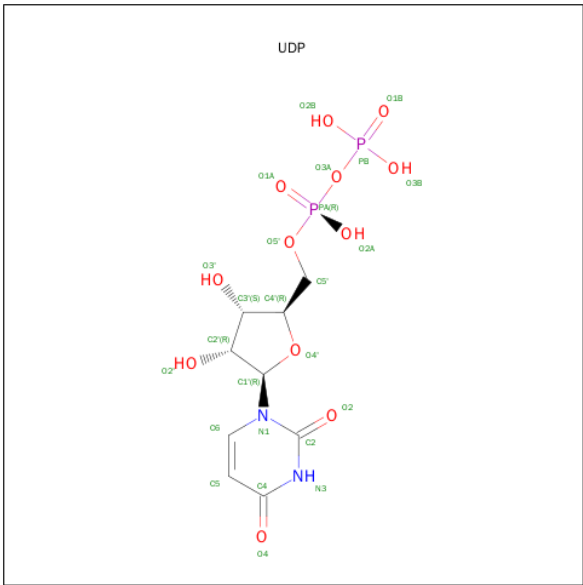
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Chain	Residue	Modelled	Actual	Comment	Reference
D	-16	SER	-	expression tag	UNP P27707
D	-15	HIS	-	expression tag	UNP P27707
D	-14	HIS	-	expression tag	UNP P27707
D	-13	HIS	-	expression tag	UNP P27707
D	-12	HIS	-	expression tag	UNP P27707
D	-11	HIS	-	expression tag	UNP P27707
D	-10	HIS	-	expression tag	UNP P27707
D	-9	SER	-	expression tag	UNP P27707
D	-8	SER	-	expression tag	UNP P27707
D	-7	GLY	-	expression tag	UNP P27707
D	-6	LEU	-	expression tag	UNP P27707
D	-5	VAL	-	expression tag	UNP P27707
D	-4	PRO	-	expression tag	UNP P27707
D	-3	ARG	-	expression tag	UNP P27707
D	-2	GLY	-	expression tag	UNP P27707
D	-1	SER	-	expression tag	UNP P27707
D	0	HIS	-	expression tag	UNP P27707

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

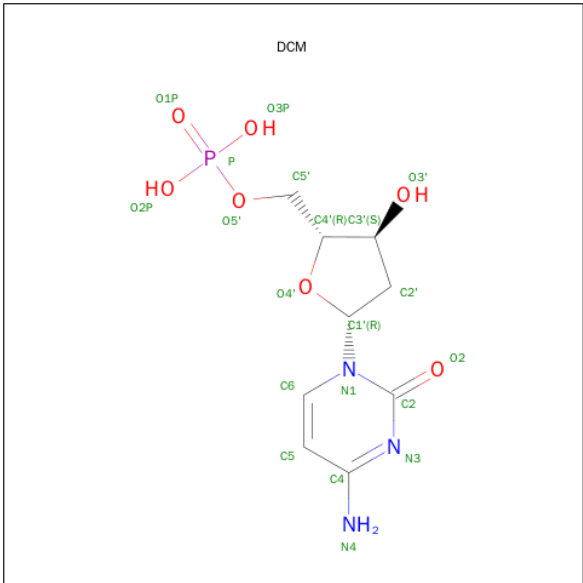
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Mg 1 1	0	0
2	A	1	Total Mg 1 1	0	0
2	D	1	Total Mg 1 1	0	0
2	C	1	Total Mg 1 1	0	0

- Molecule 3 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: C₉H₁₄N₂O₁₂P₂).

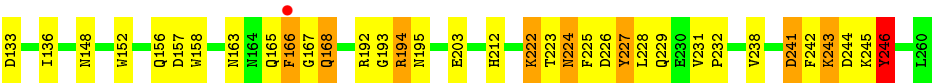


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
3	B	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
3	C	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
3	D	1	Total	C	N	O	P	0	0
			25	9	2	12	2		

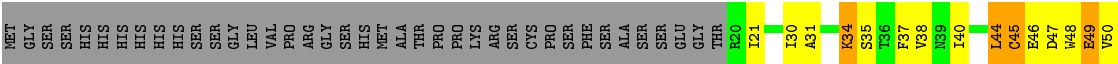
- Molecule 4 is 2'-DEOXYCYTIDINE-5'-MONOPHOSPHATE (three-letter code: DCM) (formula: C₉H₁₄N₃O₇P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			20	9	3	7	1		
4	B	1	Total	C	N	O	P	0	0
			20	9	3	7	1		
4	C	1	Total	C	N	O	P	0	0
			20	9	3	7	1		
4	D	1	Total	C	N	O	P	0	0
			20	9	3	7	1		



● Molecule 1: Deoxycytidine kinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	94.56 Å 94.56 Å 335.33 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.12 – 3.40 48.11 – 3.20	Depositor EDS
% Data completeness (in resolution range)	89.9 (48.12-3.40) 82.9 (48.11-3.20)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 3.19 Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.217 , 0.269 0.218 , 0.259	Depositor DCC
R_{free} test set	954 reflections (4.84%)	DCC
Wilson B-factor (Å ²)	65.6	Xtriage
Anisotropy	0.403	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 68.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 23589 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7808	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.97% 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: MG, DCM, UDP

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	0.41	0/1952	0.62	0/2641
1	B	0.41	0/1952	0.61	0/2641
1	C	0.42	0/1952	0.61	0/2641
1	D	0.42	0/1952	0.61	0/2641
All	All	0.41	0/7808	0.61	0/10564

There are no bond length outliers.

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	1906	0	1854	71	0
1	B	1906	0	1855	70	0
1	C	1906	0	1855	80	0
1	D	1906	0	1855	70	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	25	0	11	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	25	0	11	6	0
3	C	25	0	11	4	0
3	D	25	0	11	4	0
4	A	20	0	12	1	0
4	B	20	0	12	1	0
4	C	20	0	12	1	0
4	D	20	0	12	1	0
All	All	7808	0	7511	269	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 18.

All (269) 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:222:LYS:HE2	1:A:222:LYS:HA	1.54	0.88
1:D:222:LYS:HE2	1:D:222:LYS:HA	1.56	0.87
1:B:222:LYS:HE2	1:B:222:LYS:HA	1.56	0.84
1:C:222:LYS:HA	1:C:222:LYS:HE2	1.58	0.84
1:D:53:GLU:HG2	1:D:104:ARG:HH12	1.44	0.80
1:B:43:GLN:NE2	1:C:195:ASN:HB2	1.97	0.79
1:D:128:ARG:HG3	1:D:128:ARG:HH11	1.49	0.78
1:C:53:GLU:HG2	1:C:104:ARG:HH12	1.49	0.78
1:C:20:ARG:HD2	1:C:20:ARG:O	1.84	0.77
1:A:53:GLU:HG2	1:A:104:ARG:HH12	1.51	0.76
1:B:53:GLU:HG2	1:B:104:ARG:HH12	1.51	0.75
1:B:152:TRP:O	1:B:156:GLN:HG3	1.87	0.74
1:B:43:GLN:HE22	1:C:195:ASN:HB2	1.53	0.73
1:C:152:TRP:O	1:C:156:GLN:HG3	1.88	0.72
1:A:152:TRP:O	1:A:156:GLN:HG3	1.90	0.72
1:C:241:ASP:HA	3:C:402:UDP:O4	1.91	0.70
1:A:34:LYS:HD3	3:A:402:UDP:O2B	1.92	0.70
1:D:152:TRP:O	1:D:156:GLN:HG3	1.91	0.69
1:B:128:ARG:HG3	1:B:128:ARG:HH11	1.57	0.69
1:C:242:PHE:O	1:C:244:ASP:N	2.27	0.68
1:B:242:PHE:O	1:B:244:ASP:N	2.27	0.67
1:C:77:ASN:HD22	1:C:78:GLY:N	1.93	0.67
1:A:242:PHE:O	1:A:244:ASP:N	2.27	0.67
1:D:242:PHE:O	1:D:244:ASP:N	2.29	0.66
1:C:165:GLN:HG3	1:D:166:PHE:HE2	1.61	0.66
1:D:128:ARG:NH1	1:D:128:ARG:HG3	2.09	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:34:LYS:HD3	3:B:402:UDP:O2B	1.95	0.66
1:A:128:ARG:HH11	1:A:128:ARG:HG3	1.59	0.65
1:C:128:ARG:HH11	1:C:128:ARG:HG3	1.62	0.64
1:C:222:LYS:CE	1:C:222:LYS:HA	2.28	0.64
1:C:166:PHE:HE2	1:D:165:GLN:HG3	1.63	0.64
1:D:243:LYS:HG2	1:D:243:LYS:O	1.98	0.64
1:C:245:LYS:O	1:C:245:LYS:HG2	1.98	0.64
1:B:43:GLN:NE2	1:C:195:ASN:CB	2.60	0.63
1:D:222:LYS:HA	1:D:222:LYS:CE	2.27	0.63
1:C:243:LYS:HG2	1:C:243:LYS:O	1.97	0.63
1:B:245:LYS:HG2	1:B:245:LYS:O	1.99	0.63
1:C:242:PHE:O	1:C:246:TYR:HB3	1.99	0.63
1:C:223:THR:HG21	1:C:228:LEU:HD12	1.80	0.63
1:A:222:LYS:CE	1:A:222:LYS:HA	2.26	0.62
1:A:243:LYS:HG2	1:A:243:LYS:O	1.98	0.62
1:B:243:LYS:O	1:B:243:LYS:HG2	1.97	0.62
1:A:242:PHE:O	1:A:246:TYR:HB3	2.00	0.61
1:B:43:GLN:NE2	1:C:195:ASN:CG	2.53	0.61
1:A:128:ARG:NH1	1:A:128:ARG:HG3	2.15	0.61
1:B:242:PHE:O	1:B:246:TYR:HB3	2.00	0.61
1:D:34:LYS:N	3:D:402:UDP:O3B	2.33	0.61
1:D:245:LYS:O	1:D:245:LYS:HG2	2.01	0.61
1:A:223:THR:HG21	1:A:228:LEU:HD12	1.83	0.61
1:A:245:LYS:O	1:A:245:LYS:HG2	1.99	0.61
1:C:242:PHE:C	1:C:244:ASP:H	2.04	0.60
1:B:242:PHE:C	1:B:244:ASP:H	2.04	0.60
1:B:128:ARG:HG3	1:B:128:ARG:NH1	2.16	0.60
1:C:238:VAL:HB	1:C:242:PHE:CZ	2.36	0.60
1:D:238:VAL:HB	1:D:242:PHE:CZ	2.37	0.60
1:B:222:LYS:CE	1:B:222:LYS:HA	2.28	0.60
1:D:223:THR:HG21	1:D:228:LEU:HD12	1.83	0.59
1:B:238:VAL:HB	1:B:242:PHE:CZ	2.37	0.59
1:D:242:PHE:O	1:D:246:TYR:HB3	2.02	0.59
1:D:44:LEU:O	1:D:45:CYS:HB2	2.03	0.59
1:B:120:GLU:HG3	1:B:121:LYS:H	1.68	0.59
1:B:34:LYS:N	3:B:402:UDP:O3B	2.36	0.58
1:B:44:LEU:O	1:B:45:CYS:HB2	2.02	0.58
1:B:223:THR:HG21	1:B:228:LEU:HD12	1.85	0.58
1:D:120:GLU:HG3	1:D:121:LYS:H	1.68	0.58
1:A:242:PHE:C	1:A:244:ASP:H	2.06	0.58
1:C:120:GLU:HG3	1:C:121:LYS:H	1.68	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:120:GLU:HG3	1:A:121:LYS:H	1.68	0.58
1:C:128:ARG:HG3	1:C:128:ARG:NH1	2.18	0.58
1:A:35:SER:N	3:A:402:UDP:O3B	2.37	0.58
1:D:242:PHE:C	1:D:244:ASP:H	2.07	0.58
1:B:37:PHE:HB2	1:B:242:PHE:HE2	1.67	0.57
1:A:238:VAL:HB	1:A:242:PHE:CZ	2.39	0.57
1:C:44:LEU:O	1:C:45:CYS:HB2	2.04	0.57
1:A:44:LEU:O	1:A:45:CYS:HB2	2.05	0.56
1:A:83:GLN:HG2	1:A:87:GLU:OE2	2.06	0.56
1:A:37:PHE:HB2	1:A:242:PHE:HE2	1.70	0.56
1:D:83:GLN:HG2	1:D:87:GLU:OE2	2.07	0.55
1:C:37:PHE:HB2	1:C:242:PHE:HE2	1.71	0.55
1:D:37:PHE:HB2	1:D:242:PHE:HE2	1.72	0.55
1:C:106:ARG:NH1	1:D:157:ASP:OD2	2.39	0.54
1:B:43:GLN:HE22	1:C:195:ASN:CB	2.17	0.54
1:D:80:ASN:O	1:D:84:MET:HG3	2.07	0.54
1:C:77:ASN:C	1:C:77:ASN:HD22	2.08	0.54
1:D:34:LYS:HD3	3:D:402:UDP:O2B	2.08	0.54
1:D:241:ASP:HA	3:D:402:UDP:O4	2.08	0.54
1:C:157:ASP:OD2	1:D:106:ARG:NH1	2.41	0.54
1:D:31:ALA:HA	3:D:402:UDP:O3A	2.07	0.53
1:A:20:ARG:HD2	1:A:20:ARG:O	2.09	0.53
1:C:109:LEU:HD11	1:C:166:PHE:HB3	1.90	0.53
1:D:60:ASN:OD1	1:D:78:GLY:N	2.41	0.53
1:B:128:ARG:NH1	4:B:403:DCM:O2P	2.38	0.53
1:D:53:GLU:OE1	1:D:128:ARG:NH1	2.40	0.53
1:B:115:LYS:NZ	3:C:402:UDP:H5'2	2.24	0.53
1:D:109:LEU:HD11	1:D:166:PHE:HB3	1.91	0.53
1:C:47:ASP:OD2	1:C:119:ALA:HB1	2.09	0.53
1:D:127:GLU:O	1:D:128:ARG:HB2	2.09	0.52
1:B:20:ARG:O	1:B:20:ARG:HG2	2.08	0.52
1:A:61:VAL:HG22	1:A:79:GLY:H	1.73	0.52
1:B:47:ASP:OD2	1:B:119:ALA:HB1	2.09	0.52
1:D:53:GLU:HG2	1:D:104:ARG:NH1	2.20	0.52
1:C:83:GLN:HG2	1:C:87:GLU:OE2	2.09	0.52
1:C:166:PHE:HZ	1:D:166:PHE:HZ	1.56	0.52
1:B:193:GLY:O	1:B:194:ARG:C	2.48	0.52
1:B:83:GLN:HG2	1:B:87:GLU:OE2	2.09	0.52
1:A:109:LEU:HD11	1:A:166:PHE:HB3	1.92	0.52
1:B:90:GLU:N	1:B:90:GLU:OE1	2.39	0.51
1:A:53:GLU:OE1	1:A:128:ARG:NH1	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:53:GLU:HG2	1:B:104:ARG:NH1	2.24	0.51
1:C:242:PHE:C	1:C:244:ASP:N	2.62	0.51
1:A:47:ASP:OD2	1:A:119:ALA:HB1	2.11	0.51
1:B:109:LEU:HD11	1:B:166:PHE:HB3	1.91	0.51
1:A:148:ASN:HD21	1:B:92:TRP:HH2	1.58	0.51
1:A:242:PHE:C	1:A:244:ASP:N	2.63	0.51
1:A:120:GLU:HG3	1:A:121:LYS:N	2.26	0.51
1:A:158:TRP:HD1	1:B:102:LEU:HD22	1.75	0.51
1:A:49:GLU:HB2	1:A:116:LEU:HD22	1.93	0.51
1:C:34:LYS:H	1:C:34:LYS:CD	2.24	0.51
1:D:49:GLU:HB2	1:D:116:LEU:HD22	1.93	0.51
1:A:166:PHE:O	1:A:168:GLN:N	2.44	0.51
1:A:238:VAL:HB	1:A:242:PHE:CE1	2.47	0.50
1:B:34:LYS:CD	1:B:34:LYS:H	2.23	0.50
1:C:53:GLU:OE1	1:C:128:ARG:NH1	2.42	0.50
1:B:242:PHE:C	1:B:244:ASP:N	2.63	0.50
1:C:238:VAL:HB	1:C:242:PHE:CE1	2.46	0.50
1:A:241:ASP:HA	3:A:402:UDP:O4	2.12	0.50
1:B:238:VAL:HB	1:B:242:PHE:CE1	2.47	0.50
1:D:90:GLU:OE1	1:D:90:GLU:N	2.42	0.50
1:D:34:LYS:CD	1:D:34:LYS:H	2.24	0.50
1:B:192:ARG:HG2	1:B:193:GLY:N	2.27	0.50
1:B:53:GLU:OE1	1:B:128:ARG:NH1	2.41	0.50
1:C:166:PHE:O	1:C:168:GLN:N	2.45	0.50
1:A:34:LYS:CD	1:A:34:LYS:H	2.24	0.50
1:B:166:PHE:O	1:B:168:GLN:N	2.45	0.50
1:D:120:GLU:HG3	1:D:121:LYS:N	2.26	0.49
1:C:120:GLU:HG3	1:C:121:LYS:N	2.27	0.49
1:D:47:ASP:OD2	1:D:119:ALA:HB1	2.12	0.49
1:D:166:PHE:O	1:D:168:GLN:N	2.44	0.49
1:B:49:GLU:HB2	1:B:116:LEU:HD22	1.94	0.49
1:C:46:GLU:OE2	1:C:46:GLU:N	2.45	0.49
1:C:148:ASN:HD21	1:D:92:TRP:HH2	1.60	0.49
1:C:34:LYS:N	3:C:402:UDP:O3B	2.46	0.49
1:B:120:GLU:HG3	1:B:121:LYS:N	2.27	0.49
1:A:90:GLU:N	1:A:90:GLU:OE1	2.41	0.49
1:C:132:SER:O	1:C:136:ILE:HB	2.13	0.49
1:D:46:GLU:OE2	1:D:46:GLU:N	2.45	0.49
1:C:49:GLU:HB2	1:C:116:LEU:HD22	1.95	0.49
1:D:53:GLU:OE1	1:D:128:ARG:HG3	2.13	0.49
1:D:21:ILE:HG23	1:D:122:PRO:HB2	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:80:ASN:O	1:C:84:MET:HG3	2.13	0.48
1:C:90:GLU:OE1	1:C:90:GLU:N	2.44	0.48
1:D:193:GLY:O	1:D:194:ARG:C	2.51	0.48
1:D:238:VAL:HB	1:D:242:PHE:CE1	2.48	0.48
1:C:193:GLY:O	1:C:194:ARG:C	2.52	0.48
1:D:86:TYR:HH	4:D:403:DCM:HO3'	1.60	0.48
1:D:242:PHE:C	1:D:244:ASP:N	2.64	0.48
1:A:46:GLU:N	1:A:46:GLU:OE2	2.46	0.48
1:A:193:GLY:O	1:A:194:ARG:C	2.52	0.48
1:C:47:ASP:O	1:C:48:TRP:HD1	1.97	0.48
1:C:35:SER:O	1:C:38:VAL:HG12	2.14	0.48
1:A:34:LYS:N	3:A:402:UDP:O3B	2.47	0.47
1:C:192:ARG:HG2	1:C:193:GLY:N	2.28	0.47
1:D:132:SER:O	1:D:136:ILE:HB	2.14	0.47
1:C:158:TRP:HE1	1:D:158:TRP:HE1	1.61	0.47
1:A:128:ARG:NH1	4:A:403:DCM:O2P	2.46	0.47
1:C:166:PHE:HZ	1:D:166:PHE:CZ	2.32	0.47
1:C:166:PHE:CZ	1:D:166:PHE:CZ	3.02	0.47
1:D:192:ARG:HG2	1:D:193:GLY:N	2.29	0.47
1:B:80:ASN:O	1:B:84:MET:HG3	2.13	0.47
1:C:92:TRP:HH2	1:D:148:ASN:HD21	1.61	0.47
1:B:46:GLU:OE2	1:B:46:GLU:N	2.46	0.47
1:A:127:GLU:O	1:A:128:ARG:HB2	2.14	0.47
1:B:21:ILE:HG23	1:B:122:PRO:HB2	1.96	0.47
1:A:53:GLU:HG2	1:A:104:ARG:NH1	2.24	0.47
1:B:35:SER:O	1:B:38:VAL:HG12	2.14	0.47
1:D:128:ARG:NH1	1:D:128:ARG:CG	2.76	0.46
1:A:50:VAL:HG12	1:A:51:VAL:N	2.29	0.46
1:D:47:ASP:O	1:D:48:TRP:HD1	1.98	0.46
1:A:21:ILE:HG23	1:A:122:PRO:HB2	1.96	0.46
1:C:127:GLU:O	1:C:128:ARG:HB2	2.16	0.46
1:C:227:TYR:C	1:C:229:GLN:N	2.68	0.46
1:C:21:ILE:HG23	1:C:122:PRO:HB2	1.98	0.46
1:A:61:VAL:O	1:A:62:GLN:OE1	2.34	0.46
1:A:47:ASP:O	1:A:48:TRP:HD1	1.97	0.46
1:D:30:ILE:O	1:D:31:ALA:HB3	2.16	0.46
1:A:92:TRP:HH2	1:B:148:ASN:HD21	1.64	0.46
1:D:40:ILE:HG21	1:D:246:TYR:CD1	2.51	0.46
1:A:136:ILE:HG21	1:A:212:HIS:CE1	2.51	0.46
1:A:242:PHE:N	3:A:402:UDP:O4	2.49	0.45
1:A:35:SER:O	1:A:38:VAL:HG12	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:192:ARG:HG2	1:A:193:GLY:N	2.30	0.45
1:A:58:TRP:O	1:A:81:VAL:HG12	2.17	0.45
1:A:104:ARG:O	1:A:108:GLN:HG3	2.16	0.45
1:B:47:ASP:O	1:B:48:TRP:HD1	1.98	0.45
1:B:127:GLU:O	1:B:128:ARG:HB2	2.17	0.45
1:C:136:ILE:HG21	1:C:212:HIS:CE1	2.51	0.45
1:B:62:GLN:O	1:B:63:SER:HB2	2.17	0.45
1:C:50:VAL:HG12	1:C:51:VAL:N	2.30	0.45
1:D:231:VAL:HA	1:D:232:PRO:HD3	1.78	0.45
1:B:50:VAL:HG12	1:B:51:VAL:N	2.32	0.44
1:C:166:PHE:CE2	1:D:165:GLN:HG3	2.48	0.44
1:A:132:SER:O	1:A:136:ILE:HB	2.16	0.44
1:C:203:GLU:CD	1:C:203:GLU:H	2.21	0.44
1:C:53:GLU:HG2	1:C:104:ARG:NH1	2.24	0.44
1:A:61:VAL:O	1:A:62:GLN:HG2	2.17	0.44
1:C:31:ALA:HA	3:C:402:UDP:O3A	2.18	0.44
1:D:136:ILE:CD1	1:D:211:LYS:HB3	2.48	0.44
1:B:53:GLU:OE1	1:B:128:ARG:HG3	2.18	0.44
1:B:227:TYR:C	1:B:229:GLN:N	2.71	0.44
1:D:35:SER:O	1:D:38:VAL:HG12	2.17	0.44
1:D:104:ARG:O	1:D:108:GLN:HG3	2.18	0.43
1:A:221:LEU:C	1:A:221:LEU:HD23	2.38	0.43
1:A:227:TYR:C	1:A:229:GLN:N	2.71	0.43
1:D:50:VAL:HG12	1:D:51:VAL:N	2.32	0.43
1:B:231:VAL:HA	1:B:232:PRO:HD3	1.78	0.43
1:D:227:TYR:C	1:D:229:GLN:N	2.72	0.43
1:A:91:ARG:O	1:A:91:ARG:HD2	2.19	0.43
1:C:30:ILE:O	1:C:31:ALA:HB3	2.18	0.43
1:A:223:THR:HB	1:A:224:ASN:H	1.39	0.43
1:C:225:PHE:N	1:C:225:PHE:CD1	2.87	0.43
1:B:104:ARG:O	1:B:108:GLN:HG3	2.18	0.43
1:B:31:ALA:HA	3:B:402:UDP:O3A	2.19	0.43
1:C:104:ARG:O	1:C:108:GLN:HG3	2.19	0.43
1:B:242:PHE:N	3:B:402:UDP:O4	2.52	0.43
1:B:203:GLU:H	1:B:203:GLU:CD	2.22	0.42
1:C:128:ARG:NH1	4:C:403:DCM:O2P	2.50	0.42
1:A:110:ALA:C	1:A:112:LEU:N	2.73	0.42
1:A:231:VAL:HA	1:A:232:PRO:HD3	1.79	0.42
1:B:136:ILE:CD1	1:B:211:LYS:HB3	2.49	0.42
1:C:77:ASN:ND2	1:C:77:ASN:C	2.73	0.42
1:C:110:ALA:C	1:C:112:LEU:N	2.72	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:110:ALA:C	1:B:112:LEU:N	2.72	0.42
1:C:166:PHE:CZ	1:D:166:PHE:HZ	2.35	0.42
1:B:51:VAL:HG23	1:B:124:LEU:HD11	2.02	0.42
1:B:132:SER:O	1:B:136:ILE:HB	2.19	0.42
1:B:58:TRP:O	1:B:81:VAL:HG12	2.19	0.42
1:C:53:GLU:OE1	1:C:128:ARG:HG3	2.20	0.42
1:B:136:ILE:HG21	1:B:212:HIS:CE1	2.55	0.42
1:C:224:ASN:O	1:C:224:ASN:ND2	2.53	0.42
1:A:245:LYS:O	1:A:247:GLU:N	2.53	0.42
1:C:130:VAL:HG21	1:C:163:ASN:HD21	1.85	0.42
1:A:225:PHE:N	1:A:225:PHE:CD1	2.88	0.42
1:B:130:VAL:HG21	1:B:163:ASN:HD21	1.85	0.42
1:A:136:ILE:CD1	1:A:211:LYS:HB3	2.49	0.41
1:D:58:TRP:O	1:D:81:VAL:HG12	2.19	0.41
1:D:136:ILE:HG21	1:D:212:HIS:CE1	2.54	0.41
1:D:89:PRO:C	1:D:91:ARG:H	2.24	0.41
1:A:80:ASN:O	1:A:84:MET:HG3	2.20	0.41
1:B:241:ASP:HA	3:B:402:UDP:O4	2.20	0.41
1:B:35:SER:N	3:B:402:UDP:O3B	2.51	0.41
1:A:203:GLU:CD	1:A:203:GLU:H	2.23	0.41
1:A:112:LEU:HA	1:A:112:LEU:HD23	1.87	0.41
1:A:102:LEU:HD22	1:B:158:TRP:HD1	1.85	0.41
1:D:130:VAL:HG21	1:D:163:ASN:HD21	1.86	0.41
1:C:148:ASN:ND2	1:D:92:TRP:HH2	2.19	0.41
1:A:50:VAL:CG1	1:A:51:VAL:N	2.84	0.41
1:B:39:ASN:O	1:B:43:GLN:HG2	2.21	0.41
1:B:30:ILE:O	1:B:31:ALA:HB3	2.20	0.41
1:A:30:ILE:O	1:A:31:ALA:HB3	2.21	0.41
1:A:39:ASN:O	1:A:43:GLN:HG2	2.21	0.41
1:C:89:PRO:C	1:C:91:ARG:H	2.25	0.41
1:A:53:GLU:OE1	1:A:128:ARG:HG3	2.21	0.40
1:C:223:THR:HB	1:C:224:ASN:H	1.38	0.40
1:A:20:ARG:HD2	1:A:20:ARG:C	2.41	0.40
1:D:91:ARG:HD2	1:D:91:ARG:O	2.21	0.40
1:C:231:VAL:HA	1:C:232:PRO:HD3	1.79	0.40
1:C:117:LYS:HE2	1:C:117:LYS:HB3	1.91	0.40
1:C:20:ARG:HD2	1:C:20:ARG:C	2.41	0.40
1:C:58:TRP:O	1:C:81:VAL:HG12	2.20	0.40

There are no symmetry-related clashes.

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	224/280 (80%)	184 (82%)	32 (14%)	8 (4%)	4	34
1	B	224/280 (80%)	185 (83%)	33 (15%)	6 (3%)	6	41
1	C	224/280 (80%)	188 (84%)	30 (13%)	6 (3%)	6	41
1	D	224/280 (80%)	187 (84%)	29 (13%)	8 (4%)	4	34
All	All	896/1120 (80%)	744 (83%)	124 (14%)	28 (3%)	5	39

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	194	ARG
1	A	243	LYS
1	B	194	ARG
1	B	243	LYS
1	C	194	ARG
1	C	243	LYS
1	D	194	ARG
1	D	243	LYS
1	A	78	GLY
1	A	167	GLY
1	A	226	ASP
1	B	167	GLY
1	B	226	ASP
1	C	167	GLY
1	C	226	ASP
1	D	167	GLY
1	D	226	ASP
1	A	246	TYR
1	B	45	CYS
1	B	246	TYR
1	D	45	CYS
1	D	241	ASP

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Mol	Chain	Res	Type
1	A	45	CYS
1	C	45	CYS
1	C	246	TYR
1	D	246	TYR
1	A	195	ASN
1	D	195	ASN

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	210/256 (82%)	195 (93%)	15 (7%)	18	58
1	B	210/256 (82%)	196 (93%)	14 (7%)	20	60
1	C	210/256 (82%)	193 (92%)	17 (8%)	15	51
1	D	210/256 (82%)	195 (93%)	15 (7%)	18	58
All	All	840/1024 (82%)	779 (93%)	61 (7%)	17	56

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	ARG
1	A	34	LYS
1	A	44	LEU
1	A	49	GLU
1	A	57	ARG
1	A	112	LEU
1	A	122	PRO
1	A	133	ASP
1	A	166	PHE
1	A	168	GLN
1	A	222	LYS
1	A	224	ASN
1	A	227	TYR
1	A	241	ASP
1	A	246	TYR

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Mol	Chain	Res	Type
1	B	34	LYS
1	B	44	LEU
1	B	49	GLU
1	B	57	ARG
1	B	77	ASN
1	B	112	LEU
1	B	133	ASP
1	B	166	PHE
1	B	168	GLN
1	B	222	LYS
1	B	224	ASN
1	B	227	TYR
1	B	241	ASP
1	B	246	TYR
1	C	20	ARG
1	C	34	LYS
1	C	44	LEU
1	C	49	GLU
1	C	57	ARG
1	C	63	SER
1	C	77	ASN
1	C	112	LEU
1	C	122	PRO
1	C	133	ASP
1	C	166	PHE
1	C	168	GLN
1	C	222	LYS
1	C	224	ASN
1	C	227	TYR
1	C	241	ASP
1	C	246	TYR
1	D	34	LYS
1	D	44	LEU
1	D	49	GLU
1	D	57	ARG
1	D	63	SER
1	D	112	LEU
1	D	122	PRO
1	D	133	ASP
1	D	166	PHE
1	D	168	GLN
1	D	222	LYS

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Mol	Chain	Res	Type
1	D	224	ASN
1	D	227	TYR
1	D	241	ASP
1	D	246	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
1	A	62	GLN
1	A	77	ASN
1	A	163	ASN
1	A	198	GLN
1	A	224	ASN
1	B	43	GLN
1	B	77	ASN
1	B	163	ASN
1	B	224	ASN
1	C	43	GLN
1	C	77	ASN
1	C	163	ASN
1	C	224	ASN
1	D	43	GLN
1	D	163	ASN
1	D	198	GLN
1	D	224	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 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$
3	UDP	A	402	-	18,26,26	1.44	3 (16%)	26,40,40	2.16	4 (15%)
4	DCM	A	403	-	17,21,21	1.37	4 (23%)	24,31,31	1.20	2 (8%)
3	UDP	B	402	-	18,26,26	1.56	3 (16%)	26,40,40	2.16	4 (15%)
4	DCM	B	403	-	17,21,21	1.33	3 (17%)	24,31,31	1.18	2 (8%)
3	UDP	C	402	2	18,26,26	1.36	2 (11%)	26,40,40	2.17	4 (15%)
4	DCM	C	403	-	17,21,21	1.38	4 (23%)	24,31,31	1.18	2 (8%)
3	UDP	D	402	-	18,26,26	1.57	5 (27%)	26,40,40	2.15	4 (15%)
4	DCM	D	403	-	17,21,21	1.44	4 (23%)	24,31,31	1.18	2 (8%)

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
3	UDP	A	402	-	-	0/12/32/32	0/2/2/2
4	DCM	A	403	-	-	0/6/22/22	0/2/2/2
3	UDP	B	402	-	-	0/12/32/32	0/2/2/2
4	DCM	B	403	-	-	0/6/22/22	0/2/2/2
3	UDP	C	402	2	-	0/12/32/32	0/2/2/2
4	DCM	C	403	-	-	0/6/22/22	0/2/2/2
3	UDP	D	402	-	-	0/12/32/32	0/2/2/2
4	DCM	D	403	-	-	0/6/22/22	0/2/2/2

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	UDP	PB-O2B	-2.73	1.44	1.54
3	B	402	UDP	PB-O2B	-2.54	1.45	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	402	UDP	O5'-C5'	-2.51	1.34	1.44
4	A	403	DCM	P-O2P	-2.28	1.46	1.54
4	B	403	DCM	P-O2P	-2.27	1.46	1.54
4	D	403	DCM	P-O2P	-2.26	1.46	1.54
4	C	403	DCM	P-O2P	-2.20	1.46	1.54
3	D	402	UDP	PB-O2B	-2.10	1.47	1.54
3	D	402	UDP	C5'-C4'	-2.01	1.45	1.51
4	C	403	DCM	O4'-C1'	2.01	1.47	1.42
4	D	403	DCM	O4'-C1'	2.06	1.47	1.42
4	A	403	DCM	O4'-C1'	2.16	1.47	1.42
4	A	403	DCM	O4'-C4'	2.27	1.50	1.45
4	D	403	DCM	O4'-C4'	2.28	1.50	1.45
4	C	403	DCM	O4'-C4'	2.41	1.50	1.45
4	B	403	DCM	O4'-C4'	2.41	1.50	1.45
3	D	402	UDP	C4-N3	2.43	1.37	1.33
3	C	402	UDP	C4-N3	2.61	1.37	1.33
3	D	402	UDP	C6-N1	2.76	1.39	1.35
3	A	402	UDP	C6-N1	2.80	1.39	1.35
4	B	403	DCM	P-O3P	2.83	1.64	1.54
4	A	403	DCM	P-O3P	2.86	1.65	1.54
3	C	402	UDP	C6-N1	2.91	1.39	1.35
3	A	402	UDP	C4-N3	2.97	1.38	1.33
3	B	402	UDP	C4-N3	3.00	1.38	1.33
4	C	403	DCM	P-O3P	3.01	1.65	1.54
4	D	403	DCM	P-O3P	3.28	1.66	1.54
3	B	402	UDP	C6-N1	3.46	1.40	1.35

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	402	UDP	PA-O3A-PB	-7.62	107.11	132.67
3	B	402	UDP	PA-O3A-PB	-7.62	107.13	132.67
3	C	402	UDP	PA-O3A-PB	-7.57	107.29	132.67
3	D	402	UDP	PA-O3A-PB	-7.55	107.36	132.67
3	C	402	UDP	C4'-O4'-C1'	-3.20	106.20	109.72
3	A	402	UDP	C4'-O4'-C1'	-3.09	106.33	109.72
3	D	402	UDP	C4'-O4'-C1'	-3.02	106.40	109.72
3	B	402	UDP	C4'-O4'-C1'	-2.86	106.58	109.72
4	B	403	DCM	N4-C4-N3	2.01	120.16	116.50
4	D	403	DCM	N4-C4-N3	2.01	120.17	116.50
4	C	403	DCM	N4-C4-N3	2.02	120.18	116.50
4	A	403	DCM	N4-C4-N3	2.03	120.20	116.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	402	UDP	O3A-PA-O5'	2.22	108.83	102.94
3	C	402	UDP	O3A-PA-O5'	2.31	109.07	102.94
3	A	402	UDP	O3A-PA-O5'	2.36	109.20	102.94
3	B	402	UDP	O3A-PA-O5'	2.37	109.23	102.94
4	B	403	DCM	C2-N3-C4	3.65	120.77	115.61
4	D	403	DCM	C2-N3-C4	3.69	120.82	115.61
4	C	403	DCM	C2-N3-C4	3.71	120.84	115.61
4	A	403	DCM	C2-N3-C4	3.73	120.88	115.61
3	A	402	UDP	C4-N3-C2	6.10	120.19	114.14
3	C	402	UDP	C4-N3-C2	6.17	120.25	114.14
3	D	402	UDP	C4-N3-C2	6.18	120.26	114.14
3	B	402	UDP	C4-N3-C2	6.20	120.28	114.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 23 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	402	UDP	5	0
4	A	403	DCM	1	0
3	B	402	UDP	6	0
4	B	403	DCM	1	0
3	C	402	UDP	4	0
4	C	403	DCM	1	0
3	D	402	UDP	4	0
4	D	403	DCM	1	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	228/280 (81%)	0.11	7 (3%) 52 48	23, 63, 132, 173	0
1	B	228/280 (81%)	-0.04	2 (0%) 85 81	26, 63, 135, 190	0
1	C	228/280 (81%)	-0.17	7 (3%) 52 48	24, 59, 134, 161	0
1	D	228/280 (81%)	-0.11	3 (1%) 79 74	26, 62, 128, 193	0
All	All	912/1120 (81%)	-0.05	19 (2%) 67 61	23, 62, 134, 193	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	21	ILE	3.0
1	C	46	GLU	2.9
1	D	166	PHE	2.7
1	C	47	ASP	2.7
1	C	118	ASP	2.6
1	A	246	TYR	2.6
1	B	77	ASN	2.5
1	C	166	PHE	2.5
1	A	250	VAL	2.4
1	C	119	ALA	2.3
1	D	117	LYS	2.3
1	C	115	LYS	2.2
1	A	123	VAL	2.2
1	A	230	GLU	2.1
1	B	255	GLU	2.1
1	A	46	GLU	2.1
1	C	117	LYS	2.1
1	D	167	GLY	2.0
1	A	47	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 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
4	DCM	C	403	20/20	0.93	0.30	4.24	51,80,89,89	0
3	UDP	C	402	25/25	0.91	0.26	0.99	66,82,87,88	0
4	DCM	A	403	20/20	0.94	0.26	0.66	51,80,89,89	0
4	DCM	D	403	20/20	0.92	0.22	0.27	51,80,89,89	0
4	DCM	B	403	20/20	0.94	0.23	0.12	51,80,89,89	0
3	UDP	D	402	25/25	0.88	0.22	-0.18	66,82,87,88	0
3	UDP	B	402	25/25	0.87	0.21	-0.29	66,82,87,88	0
3	UDP	A	402	25/25	0.90	0.21	-0.72	66,82,87,88	0
2	MG	A	401	1/1	0.96	0.12	-	56,56,56,56	0
2	MG	B	401	1/1	0.94	0.26	-	56,56,56,56	0
2	MG	C	401	1/1	0.95	0.38	-	56,56,56,56	0
2	MG	D	401	1/1	0.92	0.32	-	56,56,56,56	0

6.5 Other polymers [i](#)

There are no such residues in this entry.