



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 03:11 pm GMT

PDB ID : 2ZI7  
Title : C4S dCK variant of dCK in complex with D-dG+UDP  
Authors : Sabini, E.; Lavie, A.  
Deposited on : 2008-02-13  
Resolution : 1.97 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](mailto: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.

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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

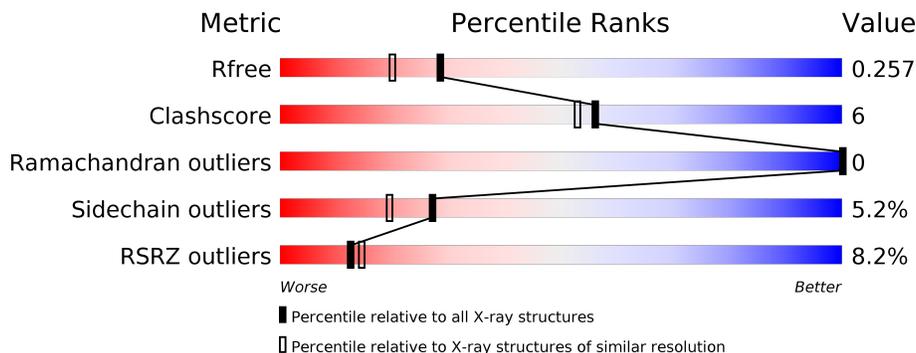
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.97 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	9293 (2.00-1.96)
Clashscore	112137	10621 (2.00-1.96)
Ramachandran outliers	110173	10502 (2.00-1.96)
Sidechain outliers	110143	10501 (2.00-1.96)
RSRZ outliers	101464	9395 (2.00-1.96)

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  $\geq 3$ , 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	279	 5% 65% 11% • 22%
1	B	279	 8% 60% 14% • 25%

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3697 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
			Total	C	N	O	S			
1	B	208	1732	1114	287	325	6	0	0	0
1	A	218	1793	1153	294	340	6	0	0	0

There are 46 discrepancies between the modelled and reference sequences:

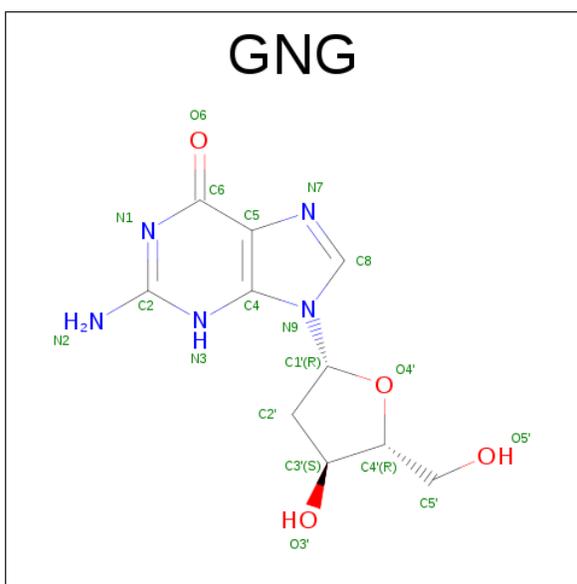
Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	MET	-	EXPRESSION TAG	UNP P27707
B	-17	GLY	-	EXPRESSION TAG	UNP P27707
B	-16	SER	-	EXPRESSION TAG	UNP P27707
B	-15	SER	-	EXPRESSION TAG	UNP P27707
B	-14	HIS	-	EXPRESSION TAG	UNP P27707
B	-13	HIS	-	EXPRESSION TAG	UNP P27707
B	-12	HIS	-	EXPRESSION TAG	UNP P27707
B	-11	HIS	-	EXPRESSION TAG	UNP P27707
B	-10	HIS	-	EXPRESSION TAG	UNP P27707
B	-9	HIS	-	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
B	9	SER	CYS	ENGINEERED	UNP P27707
B	45	SER	CYS	ENGINEERED	UNP P27707
B	59	SER	CYS	ENGINEERED	UNP P27707
B	146	SER	CYS	ENGINEERED	UNP P27707
A	-18	MET	-	EXPRESSION TAG	UNP P27707
A	-17	GLY	-	EXPRESSION TAG	UNP P27707

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-16	SER	-	EXPRESSION TAG	UNP P27707
A	-15	SER	-	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	HIS	-	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
A	9	SER	CYS	ENGINEERED	UNP P27707
A	45	SER	CYS	ENGINEERED	UNP P27707
A	59	SER	CYS	ENGINEERED	UNP P27707
A	146	SER	CYS	ENGINEERED	UNP P27707

- Molecule 2 is 2'-DEOXY-GUANOSINE (three-letter code: GNG) (formula: C<sub>10</sub>H<sub>13</sub>N<sub>5</sub>O<sub>4</sub>).



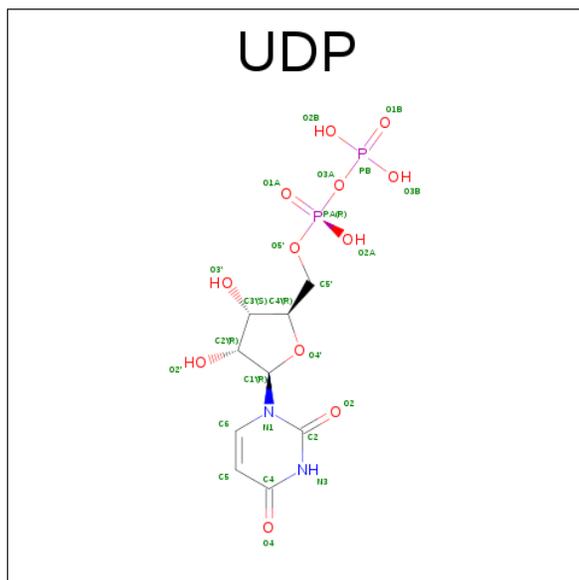
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	B	1	19	10	5	4	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	19	10	5	4	0	0

- Molecule 3 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula:  $C_9H_{14}N_2O_{12}P_2$ ).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	43	Total O 43 43	0	0
4	B	41	Total O 41 41	0	0



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	43.80Å 107.50Å 110.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 1.97 28.92 – 1.97	Depositor EDS
% Data completeness (in resolution range)	96.0 (30.00-1.97) 96.0 (28.92-1.97)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.20 (at 1.96Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.206 , 0.259 0.207 , 0.257	Depositor DCC
$R_{free}$ test set	3607 reflections (11.06%)	DCC
Wilson B-factor (Å <sup>2</sup> )	25.5	Xtrriage
Anisotropy	0.398	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.42 , 59.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.42$ , $\langle L^2 \rangle = 0.25$	Xtrriage
Estimated twinning fraction	0.038 for -h,l,k	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3697	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.70% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: UDP, GNG

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	0.81	0/1836	0.78	0/2487
1	B	0.74	1/1774 (0.1%)	0.77	0/2401
All	All	0.78	1/3610 (0.0%)	0.77	0/4888

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	183	GLU	CG-CD	6.55	1.61	1.51

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	1793	0	1728	18	0
1	B	1732	0	1673	24	0
2	A	19	0	13	1	0
2	B	19	0	13	2	0
3	A	25	0	11	0	0
3	B	25	0	11	0	0
4	A	43	0	0	0	0
4	B	41	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3697	0	3449	40	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (40) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:38:VAL:HG22	1:B:50:VAL:HG11	1.71	0.73
1:B:140:ASN:HD21	1:B:207:LYS:HD3	1.51	0.72
1:B:38:VAL:HG22	1:B:50:VAL:CG1	2.25	0.66
1:A:97:GLN:NE2	1:A:133:ASP:OD1	2.28	0.64
1:B:140:ASN:HD21	1:B:207:LYS:CD	2.13	0.60
1:A:257:LEU:HA	1:A:260:LEU:HD12	1.85	0.58
1:A:44:LEU:HD21	1:A:250:VAL:HG21	1.87	0.57
1:B:217:LEU:HD22	1:B:218:HIS:CE1	2.39	0.56
1:B:161:TRP:HB2	1:A:106:ARG:NH1	2.24	0.52
1:B:49:GLU:HG3	1:B:50:VAL:N	2.25	0.52
1:A:215:TRP:HA	1:A:221:LEU:HB3	1.93	0.51
1:A:123:VAL:HG21	1:A:257:LEU:HD13	1.94	0.50
1:A:141:LEU:HD11	2:A:502:GNG:HN22	1.76	0.50
1:B:104:ARG:NH2	2:B:503:GNG:O6	2.45	0.50
1:A:163:ASN:O	1:A:167:GLY:HA3	2.13	0.49
1:A:215:TRP:HB2	1:A:221:LEU:HD22	1.93	0.48
1:A:49:GLU:HG3	1:A:116:LEU:HD13	1.97	0.47
1:A:108:GLN:HB3	1:A:170:LEU:HD22	1.97	0.47
1:A:40:ILE:O	1:A:44:LEU:HD13	2.14	0.46
1:A:231:VAL:O	1:A:231:VAL:HG13	2.15	0.46
1:B:22:LYS:HA	1:B:173:ASP:OD2	2.15	0.46
1:B:136:ILE:HD11	1:B:215:TRP:HE3	1.81	0.46
1:B:162:MET:HE1	1:A:161:TRP:CH2	2.50	0.45
1:B:41:LEU:HA	1:B:44:LEU:HD12	1.98	0.45
1:B:26:ILE:HD11	1:B:125:PHE:HB3	1.98	0.45
1:B:57:ARG:HD2	1:B:99:TYR:OH	2.17	0.45
1:B:53:GLU:OE1	2:B:503:GNG:H8	2.17	0.44
1:B:85:MET:HG3	1:B:93:SER:HA	1.97	0.44
1:A:119:ALA:HB3	1:A:122:PRO:HB3	1.99	0.44
1:B:101:CYS:HG	1:B:158:TRP:HZ3	1.64	0.44
1:A:177:TYR:HE2	1:A:217:LEU:HD22	1.83	0.43
1:B:40:ILE:O	1:B:44:LEU:HG	2.20	0.42
1:B:49:GLU:HG3	1:B:50:VAL:H	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:SER:OG	1:A:141:LEU:HD22	2.20	0.41
1:B:102:LEU:O	1:B:106:ARG:HG3	2.20	0.41
1:B:59:SER:HA	1:B:80:ASN:OD1	2.20	0.41
1:B:179:GLN:HE22	1:B:217:LEU:HD12	1.86	0.41
1:B:180:ALA:O	4:B:541:HOH:O	2.22	0.40
1:A:247:GLU:O	1:A:251:GLU:HG3	2.21	0.40
1:B:152:TRP:O	1:B:156:GLN:HG3	2.21	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	212/279 (76%)	208 (98%)	4 (2%)	0	100	100
1	B	200/279 (72%)	193 (96%)	7 (4%)	0	100	100
All	All	412/558 (74%)	401 (97%)	11 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	195/255 (76%)	182 (93%)	13 (7%)	19	12

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	190/255 (74%)	183 (96%)	7 (4%)	39	33
All	All	385/510 (76%)	365 (95%)	20 (5%)	27	19

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	57	ARG
1	B	59	SER
1	B	111	SER
1	B	113	ASN
1	B	214	SER
1	B	252	LYS
1	B	260	LEU
1	A	38	VAL
1	A	46	GLU
1	A	62	GLN
1	A	93	SER
1	A	116	LEU
1	A	117	LYS
1	A	157	ASP
1	A	169	SER
1	A	183	GLU
1	A	217	LEU
1	A	247	GLU
1	A	252	LYS
1	A	255	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	97	GLN
1	B	140	ASN
1	B	179	GLN
1	A	60	ASN
1	A	113	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 [i](#)

4 ligands 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$
3	UDP	A	301	-	21,26,26	1.44	5 (23%)	22,40,40	1.36	5 (22%)
2	GNG	A	502	1	17,21,21	1.73	4 (23%)	20,31,31	3.46	11 (55%)
3	UDP	B	301	-	21,26,26	1.68	5 (23%)	22,40,40	1.61	3 (13%)
2	GNG	B	503	1	17,21,21	2.14	4 (23%)	20,31,31	2.82	8 (40%)

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	301	-	-	0/12/32/32	0/2/2/2
2	GNG	A	502	1	-	0/2/18/18	0/3/3/3
3	UDP	B	301	-	-	0/12/32/32	0/2/2/2
2	GNG	B	503	1	-	0/2/18/18	0/3/3/3

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	503	GNG	C5-C4	-3.06	1.33	1.40
2	A	502	GNG	C8-N7	-2.39	1.30	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	301	UDP	C2-N3	-2.35	1.33	1.38
3	A	301	UDP	PB-O3B	-2.18	1.45	1.54
3	A	301	UDP	O2'-C2'	-2.17	1.38	1.43
3	B	301	UDP	PB-O3B	-2.06	1.46	1.54
3	B	301	UDP	PB-O1B	-2.03	1.43	1.50
3	A	301	UDP	PA-O1A	-2.01	1.43	1.50
2	A	502	GNG	C4-N3	2.01	1.39	1.35
3	B	301	UDP	C4-N3	2.03	1.36	1.33
2	A	502	GNG	C2-N1	2.09	1.39	1.35
3	A	301	UDP	C4-N3	2.47	1.37	1.33
3	A	301	UDP	PB-O3A	2.65	1.64	1.60
2	B	503	GNG	O4'-C4'	3.08	1.52	1.45
2	B	503	GNG	C4-N3	3.26	1.41	1.35
3	B	301	UDP	PB-O3A	4.76	1.67	1.60
2	A	502	GNG	C6-N1	5.31	1.42	1.33
2	B	503	GNG	C6-N1	6.05	1.44	1.33

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	502	GNG	C5-C6-N1	-8.70	111.10	123.48
2	B	503	GNG	C5-C6-N1	-6.28	114.54	123.48
2	A	502	GNG	C2-N3-C4	-4.53	109.87	115.16
2	B	503	GNG	N2-C2-N1	-3.60	111.48	117.24
2	A	502	GNG	N2-C2-N1	-3.57	111.53	117.24
2	B	503	GNG	N3-C2-N1	-3.07	122.97	127.46
2	B	503	GNG	C6-C5-C4	-2.83	118.03	120.84
3	A	301	UDP	O3'-C3'-C2'	-2.38	104.20	111.83
2	A	502	GNG	N3-C2-N1	-2.34	124.04	127.46
3	A	301	UDP	O3'-C3'-C4'	-2.22	104.59	111.09
3	B	301	UDP	O3'-C3'-C4'	-2.20	104.67	111.09
2	A	502	GNG	O4'-C1'-N9	-2.02	104.37	107.78
3	B	301	UDP	O3B-PB-O1B	2.18	119.04	110.50
3	A	301	UDP	O4'-C1'-N1	2.36	112.81	108.08
3	A	301	UDP	O5'-PA-O1A	2.80	120.52	109.25
2	A	502	GNG	C2'-C3'-C4'	2.85	108.80	102.73
2	B	503	GNG	O4'-C1'-N9	2.85	112.58	107.78
2	A	502	GNG	O4'-C1'-C2'	3.12	112.23	106.25
3	A	301	UDP	C4-N3-C2	3.25	116.92	114.13
2	A	502	GNG	N2-C2-N3	3.61	124.42	117.75
2	A	502	GNG	C5'-C4'-C3'	4.17	125.42	114.84
2	B	503	GNG	N2-C2-N3	4.20	125.51	117.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	502	GNG	C4-C5-N7	4.62	113.87	109.41
2	B	503	GNG	C6-N1-C2	4.73	122.86	116.06
2	B	503	GNG	C4-C5-N7	5.21	114.44	109.41
3	B	301	UDP	C4-N3-C2	5.33	118.71	114.13
2	A	502	GNG	C6-N1-C2	6.22	125.00	116.06

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	502	GNG	1	0
2	B	503	GNG	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	218/279 (78%)	0.54	14 (6%) 20 23	13, 25, 43, 58	0
1	B	208/279 (74%)	0.65	21 (10%) 8 9	15, 28, 49, 56	0
All	All	426/558 (76%)	0.59	35 (8%) 12 14	13, 27, 48, 58	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	64	THR	10.0
1	A	220	THR	8.5
1	B	260	LEU	6.6
1	A	63	SER	5.7
1	B	77	ASN	4.8
1	B	44	LEU	4.4
1	B	220	THR	3.4
1	B	78	GLY	3.4
1	A	62	GLN	3.2
1	B	258	SER	3.2
1	A	260	LEU	3.1
1	B	217	LEU	3.0
1	A	219	ARG	3.0
1	A	221	LEU	3.0
1	B	46	GLU	2.9
1	B	244	ASP	2.8
1	B	259	THR	2.8
1	B	43	GLN	2.8
1	A	26	ILE	2.7
1	A	100	ALA	2.7
1	B	112	LEU	2.5
1	B	47	ASP	2.5
1	B	178	LEU	2.5
1	B	26	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	78	GLY	2.4
1	A	130	VAL	2.3
1	A	129	SER	2.3
1	B	219	ARG	2.3
1	B	28	GLY	2.2
1	A	116	LEU	2.2
1	B	176	ILE	2.1
1	B	25	SER	2.1
1	B	245	LYS	2.0
1	B	247	GLU	2.0
1	A	240	GLU	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q < 0.9
2	GNG	A	502	19/19	0.97	0.17	0.55	16,17,20,23	0
2	GNG	B	503	19/19	0.97	0.14	0.13	13,15,18,19	0
3	UDP	A	301	25/25	0.98	0.11	-0.74	11,20,29,30	0
3	UDP	B	301	25/25	0.98	0.11	-0.78	14,22,27,28	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.