



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 11:41 PM GMT

PDB ID : 1YA6  
Title : alpha-glucosyltransferase in complex with UDP and a 13-mer DNA containing a central A:G mismatch  
Authors : Lariviere, L.; Sommer, N.; Morera, S.  
Deposited on : 2004-12-17  
Resolution : 2.40 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (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

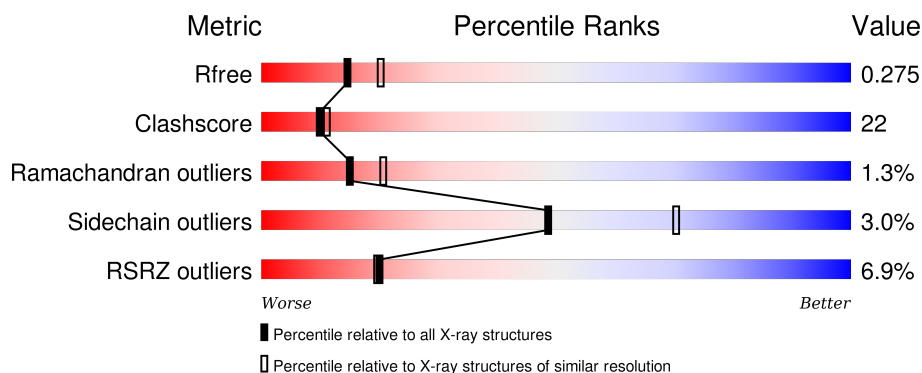
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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	C	12	<div> <div>50%</div> <div>42%</div> <div>8%</div> </div>
2	D	12	<div> <div>75%</div> <div>25%</div> </div>
3	A	403	<div> <div>7%</div> <div>60%</div> <div>32%</div> <div>• 6%</div> </div>
3	B	403	<div> <div>6%</div> <div>51%</div> <div>40%</div> <div>• 6%</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	NCO	B	3002	-	-	-	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6977 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 5'-D(\*AP\*TP\*AP\*CP\*TP\*AP\*AP\*GP\*AP\*TP\*AP\*G)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	12	Total	C	N	O	P	0	0	0
			246	119	49	67	11			

- Molecule 2 is a DNA chain called 5'-D(\*CP\*TP\*AP\*TP\*CP\*TP\*GP\*AP\*GP\*TP\*AP\*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	12	Total	C	N	O	P	0	0	0
			242	118	41	72	11			

- Molecule 3 is a protein called DNA alpha-glucosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	380	Total	C	N	O	S	0	0	0
			3118	1986	532	584	16			
3	B	377	Total	C	N	O	S	0	0	0
			3091	1970	527	577	17			

There are 6 discrepancies between the modelled and reference sequences:

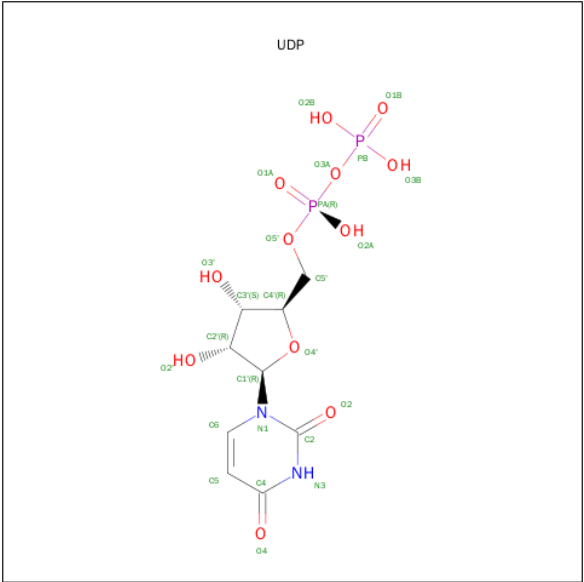
Chain	Residue	Modelled	Actual	Comment	Reference
A	998	MET	-	CLONING ARTIFACT	UNP P04519
A	999	GLY	-	CLONING ARTIFACT	UNP P04519
A	1000	SER	-	CLONING ARTIFACT	UNP P04519
B	998	MET	-	CLONING ARTIFACT	UNP P04519
B	999	GLY	-	CLONING ARTIFACT	UNP P04519
B	1000	SER	-	CLONING ARTIFACT	UNP P04519

- Molecule 4 is COBALT HEXAMMINE(III) (three-letter code: NCO) (formula: CoH<sub>18</sub>N<sub>6</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	Co	N	0	0
			7	1	6		
4	B	1	Total	Co	N	0	0
			7	1	6		
4	A	1	Total	Co	N	0	0
			7	1	6		
4	B	1	Total	Co	N	0	0
			7	1	6		
4	D	1	Total	Co	N	0	0
			7	1	6		
4	B	1	Total	Co	N	0	0
			7	1	6		
4	A	1	Total	Co	N	0	0
			7	1	6		

- Molecule 5 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: C<sub>9</sub>H<sub>14</sub>N<sub>2</sub>O<sub>12</sub>P<sub>2</sub>).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	89	Total	O	0	0
			89	89		
6	B	81	Total	O	0	0
			81	81		
6	C	5	Total	O	0	0
			5	5		
6	D	6	Total	O	0	0
			6	6		

### 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 errors 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.

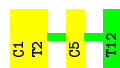
- Molecule 1: 5'-D(\*AP\*TP\*AP\*CP\*TP\*AP\*AP\*GP\*AP\*TP\*AP\*G)-3'

Chain C:



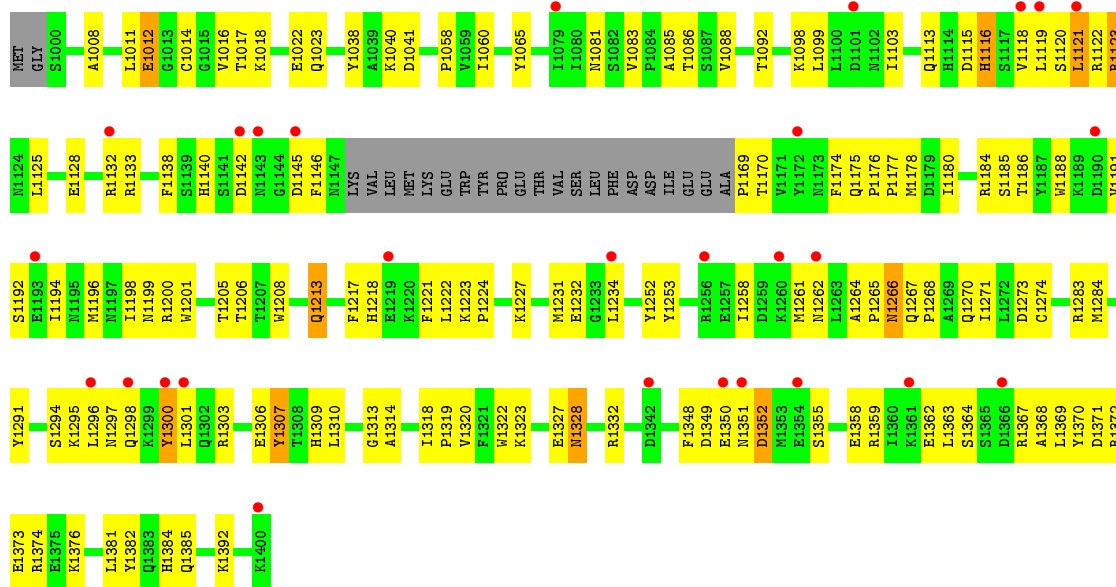
- Molecule 2: 5'-D(\*CP\*TP\*AP\*TP\*CP\*TP\*GP\*AP\*GP\*TP\*AP\*T)-3'

Chain D:



- Molecule 3: DNA alpha-glucosyltransferase

Chain A:



- Molecule 3: DNA alpha-glucosyltransferase

Chain B:

F1331	M1261	Y1096	M998
R1332	M1262	K1097	G999
V1333	Q1267	K1098	I1003
D1334	P1268	L1099	C1004
N1335	A1269	L1100	I1005
T1336	Q1270	D1101	A1008
P1337	I1271	R1108	R1009
L1338	L1272	H1114	E1012
T1339	D1273	D1115	G1013
S1340	C1274	H1116	C1014
H1341	Y1275	L1119	K1018
D1342	I1276	R1122	Q1023
I1346	E1279	R1123	R1024
W1347	M1280	N1124	D1025
D1352	R1283	L1125	W1026
M1353	M1284	G1126	K1029
E1354	F1289	E1128	V1034
S1355	G1290	E1129	Y1037
E1358	I1291	T1130	Y1038
R1359	Q1292	V1131	A1039
I1360	K1295	R1132	K1040
K1361	L1296	R1133	D1041
E1362	L1297	F1138	R1046
L1363	M1297	S1141	S1054
D1366	Q1298	ASN	V1059
R1367	K1299	GLY	I1060
A1368	Y1300	ASP	L1061
L1369	L1301	PHE	A1062
Y1370	Q1302	ASN	R1063
D1371	R1303	LYS	E1064
R1372	S1304	VAL	Y1065
E1373	L1305	LEU	D1066
K1376	E1306	MET	K1067
A1377	Y1307	GLU	A1068
Y1378	T1308	TRP	L1069
Y1382	L1310	TYR	M1081
Q1383	E1311	PRO	S1082
H1384	L1312	GLU	V1083
Q1385	G1315	THR	T1086
D1386	G1316	ILE	Q1089
S1387	T1317	GLU	E1090
C1390	I1318	ASP	A1091
F1391	P1319	ASP	T1092
V1320	V1320	ILE	I1093
F1321	F1321	GLU	M1094
W1322	W1322	ALA	M1095
K1323	K1323	P1169	
S1324	S1324	T1170	
T1325	T1325		
G1326	G1326		
E1327	E1327		
M1328	M1328		
L1329	L1329		
K1330	K1330		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	45.85Å 119.32Å 86.81Å 90.00° 92.87° 90.00°	Depositor
Resolution (Å)	20.00 – 2.40 19.98 – 2.40	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.40) 99.9 (19.98-2.40)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.30 (at 2.41Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.230 , 0.280 0.227 , 0.275	Depositor DCC
$R_{free}$ test set	1795 reflections (5.21%)	DCC
Wilson B-factor (Å <sup>2</sup> )	50.4	Xtriage
Anisotropy	0.718	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 50.6	EDS
Estimated twinning fraction	0.030 for h,-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 36229 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6977	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.49% 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, 0.375 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: UDP, NCO

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	C	0.48	0/277	0.76	0/426
2	D	0.44	0/270	0.71	0/415
3	A	0.41	0/3190	0.63	0/4301
3	B	0.39	0/3162	0.62	0/4262
All	All	0.41	0/6899	0.64	0/9404

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1
3	A	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	A	1307	TYR	Sidechain
1	C	6	DT	Sidechain

## 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	C	246	0	137	4	0
2	D	242	0	139	6	0
3	A	3118	0	3060	143	0
3	B	3091	0	3044	141	0
4	A	21	0	0	3	0
4	B	21	0	0	4	0
4	D	7	0	0	0	0
5	A	25	0	11	0	0
5	B	25	0	11	1	0
6	A	89	0	0	3	0
6	B	81	0	0	3	0
6	C	5	0	0	0	0
6	D	6	0	0	0	0
All	All	6977	0	6402	290	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 22.

The worst 5 of 290 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)
3:A:1213:GLN:H	3:A:1213:GLN:NE2	1.68	0.92
3:A:1118:VAL:HG12	3:A:1145:ASP:HB2	1.55	0.87
3:A:1113:GLN:HE22	3:A:1115:ASP:HB2	1.40	0.85
3:A:1116:HIS:CE1	4:A:2002:NCO:N3	2.45	0.85
3:A:1318:ILE:HG12	3:A:1370:TYR:CE1	2.15	0.82

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	
3	A	376/403 (93%)	336 (89%)	36 (10%)	4 (1%)	17	25
3	B	373/403 (93%)	336 (90%)	31 (8%)	6 (2%)	12	16
All	All	749/806 (93%)	672 (90%)	67 (9%)	10 (1%)	15	21

5 of 10 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	1352	ASP
3	B	1191	VAL
3	B	1335	ASN
3	A	1185	SER
3	B	1273	ASP

### 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	
3	A	347/368 (94%)	338 (97%)	9 (3%)	54	74
3	B	344/368 (94%)	332 (96%)	12 (4%)	43	64
All	All	691/736 (94%)	670 (97%)	21 (3%)	48	70

5 of 21 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	B	1012	GLU
3	B	1119	LEU
3	B	1291	TYR
3	A	1392	LYS
3	B	1292	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 37 such

sidechains are listed below:

Mol	Chain	Res	Type
3	A	1384	HIS
3	B	1094	ASN
3	B	1383	GLN
3	A	1394	GLN
3	B	1023	GLN

### 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](#)

9 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	$\# Z  > 2$
5	UDP	A	2001	-	18,26,26	1.06	1 (5%)	26,40,40	2.77	3 (11%)
4	NCO	A	2002	-	6,6,6	1.14	0	0,15,15	0.00	-
4	NCO	A	2003	-	6,6,6	1.35	0	0,15,15	0.00	-
4	NCO	A	2004	-	6,6,6	1.28	0	0,15,15	0.00	-
5	UDP	B	3001	-	18,26,26	1.07	2 (11%)	26,40,40	2.71	3 (11%)
4	NCO	B	3002	-	6,6,6	1.22	0	0,15,15	0.00	-
4	NCO	B	3003	-	6,6,6	1.29	0	0,15,15	0.00	-
4	NCO	B	3004	-	6,6,6	1.30	0	0,15,15	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	NCO	D	1003	-	6,6,6	1.39	0	0,15,15	0.00	-

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	UDP	A	2001	-	-	0/12/32/32	0/2/2/2
4	NCO	A	2002	-	-	0/0/0/0	0/0/0/0
4	NCO	A	2003	-	-	0/0/0/0	0/0/0/0
4	NCO	A	2004	-	-	0/0/0/0	0/0/0/0
5	UDP	B	3001	-	-	0/12/32/32	0/2/2/2
4	NCO	B	3002	-	-	0/0/0/0	0/0/0/0
4	NCO	B	3003	-	-	0/0/0/0	0/0/0/0
4	NCO	B	3004	-	-	0/0/0/0	0/0/0/0
4	NCO	D	1003	-	-	0/0/0/0	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	3001	UDP	C6-C5	-2.02	1.33	1.38
5	A	2001	UDP	C4-N3	2.67	1.38	1.33
5	B	3001	UDP	C4-N3	2.67	1.38	1.33

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	2001	UDP	C5-C4-N3	-3.40	114.39	123.12
5	B	3001	UDP	C5-C4-N3	-3.29	114.68	123.12
5	B	3001	UDP	O2B-PB-O1B	2.68	119.22	110.58
5	A	2001	UDP	O2B-PB-O1B	2.76	119.46	110.58
5	B	3001	UDP	C4-N3-C2	12.76	126.78	114.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	2002	NCO	3	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	3001	UDP	1	0
4	B	3002	NCO	4	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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	C	12/12 (100%)	0.22	0 <span>100</span> <span>100</span>	38, 62, 71, 72	0
2	D	12/12 (100%)	-0.18	0 <span>100</span> <span>100</span>	50, 60, 74, 74	0
3	A	380/403 (94%)	0.34	28 (7%) <span>17</span> <span>17</span>	34, 60, 87, 100	0
3	B	377/403 (93%)	0.39	26 (6%) <span>20</span> <span>19</span>	39, 65, 84, 97	0
All	All	781/830 (94%)	0.36	54 (6%) <span>20</span> <span>19</span>	34, 62, 86, 100	0

The worst 5 of 54 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	A	1143	ASN	7.3
3	A	1145	ASP	5.5
3	B	1300	TYR	4.5
3	A	1296	LEU	4.0
3	B	1132	ARG	3.6

### 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( $\text{\AA}^2$ )	Q<0.9
4	NCO	B	3002	7/7	0.93	0.22	3.63	78,79,79,79	0
4	NCO	A	2002	7/7	0.96	0.19	0.32	63,63,65,66	0
4	NCO	A	2004	7/7	0.93	0.19	0.19	92,92,93,93	7
4	NCO	B	3003	7/7	0.94	0.17	-0.47	98,99,99,99	0
4	NCO	D	1003	7/7	0.97	0.14	-0.49	88,89,89,91	0
4	NCO	A	2003	7/7	0.96	0.15	-0.51	81,82,84,84	0
5	UDP	B	3001	25/25	0.98	0.10	-1.24	37,43,47,49	0
5	UDP	A	2001	25/25	0.98	0.08	-1.40	38,44,46,48	0
4	NCO	B	3004	7/7	0.92	0.20	-	71,71,71,73	7

## 6.5 Other polymers [i](#)

There are no such residues in this entry.