



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 4, 2026 – 09:51 PM EDT

PDB ID : 1Y8Z / pdb_00001y8z
Title : alpha-glucosyltransferase in complex with UDP and a 13-mer DNA containing a HMU base at 1.9 Å resolution
Authors : Lariviere, L.; Sommer, N.; Morera, S.
Deposited on : 2004-12-14
Resolution : 1.90 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

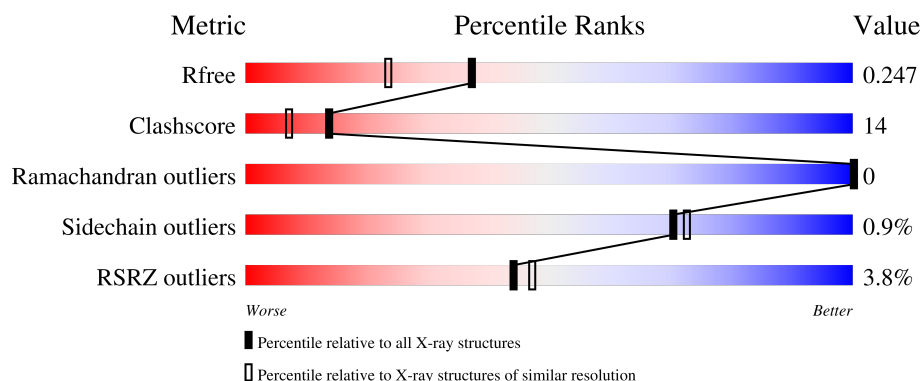
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	180053	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	13	8% 54% 46%
2	D	9	44% 56%
3	A	402	3% 70% 26% ••
3	B	402	4% 63% 27% • 7%

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
6	GOL	A	5001	-	X	-	-
6	GOL	A	5002	-	X	-	-
6	GOL	A	5003	-	X	-	-
6	GOL	B	5004	-	X	-	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 7340 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(*GP*AP*TP*AP*CP*TP*(5HU)P*AP*GP*AP*T P*AP*G)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	13	Total	C	N	O	P	0	0	0
			268	129	51	76	12			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	9	Total	C	N	O	P	0	0	0
			181	88	32	53	8			

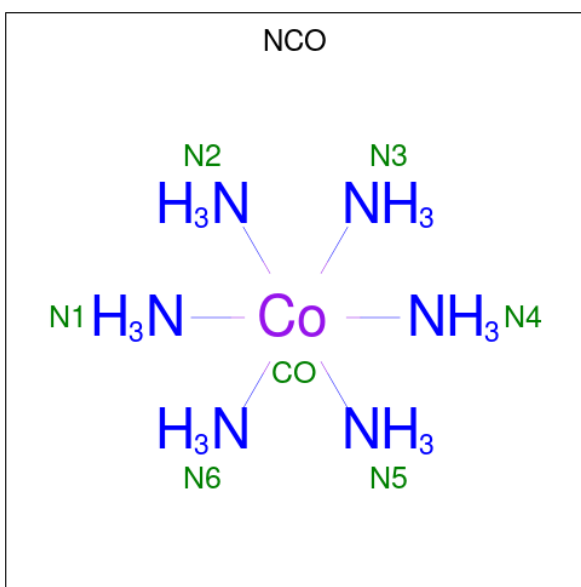
- Molecule 3 is a protein called DNA alpha-glucosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	391	Total	C	N	O	S	0	0	0
			3218	2053	546	600	19			
3	B	373	Total	C	N	O	S	0	0	0
			3069	1956	523	572	18			

There are 8 discrepancies between the modelled and reference sequences:

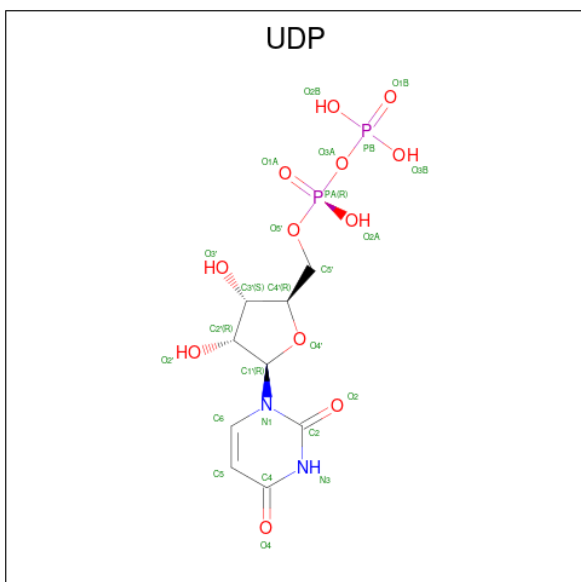
Chain	Residue	Modelled	Actual	Comment	Reference
A	999	GLY	-	cloning artifact	UNP P04519
A	1000	SER	-	cloning artifact	UNP P04519
A	1014	CME	CYS	modified residue	UNP P04519
A	1274	CME	CYS	modified residue	UNP P04519
B	999	GLY	-	cloning artifact	UNP P04519
B	1000	SER	-	cloning artifact	UNP P04519
B	1014	CME	CYS	modified residue	UNP P04519
B	1274	CME	CYS	modified residue	UNP P04519

- Molecule 4 is COBALT HEXAMMINE(III) (CCD ID: NCO) (formula: CoH₁₈N₆).



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

- Molecule 5 is URIDINE-5'-DIPHOSPHATE (CCD ID: UDP) (formula: $C_9H_{14}N_2O_{12}P_2$).



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 GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		
6	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 7 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	1	Total	Cl	0	0
			1	1		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	C	21	Total 21	O 21	0	0
8	D	8	Total 8	O 8	0	0
8	A	266	Total 266	O 266	0	0
8	B	199	Total 199	O 199	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

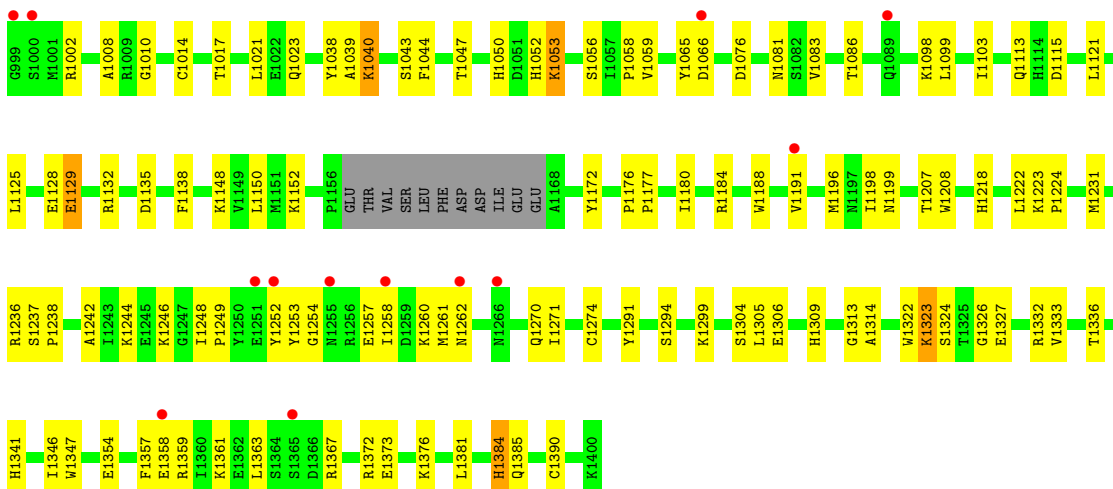
- Molecule 1: 5'-D(*GP*AP*TP*AP*CP*TP*(5HU)P*AP*GP*AP*TP*AP*G)-3'



- Molecule 2: 5'-D(*CP*TP*AP*TP*CP*TP*GP*AP*G)-3'

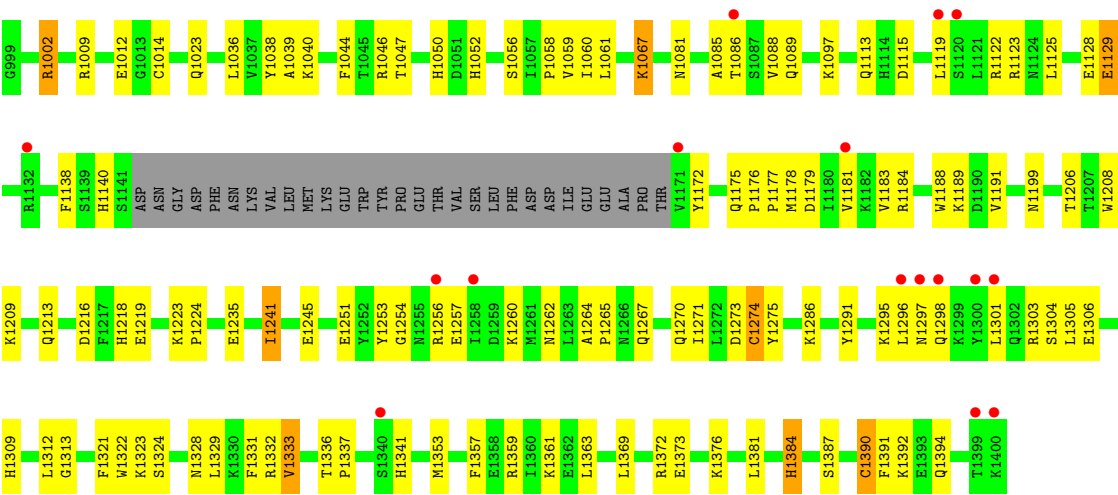


- Molecule 3: DNA alpha-glucosyltransferase



- Molecule 3: DNA alpha-glucosyltransferase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	41.26Å 114.66Å 86.14Å 90.00° 89.49° 90.00°	Depositor
Resolution (Å)	20.00 – 1.90 20.00 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.0 (20.00-1.90) 99.0 (20.00-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.45 (at 1.90Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.208 , 0.250 0.204 , 0.247	Depositor DCC
R_{free} test set	3141 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	24.9	Xtriage
Anisotropy	0.247	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 52.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.026 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7340	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.20% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 5HU, CME, NCO, GOL, CL, 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	C	0.25	0/277	0.60	0/424
2	D	0.25	0/202	0.66	0/310
3	A	0.38	0/3272	0.90	10/4410 (0.2%)
3	B	0.35	0/3117	0.87	9/4198 (0.2%)
All	All	0.36	0/6868	0.87	19/9342 (0.2%)

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1304	SER	N-CA-C	9.20	122.93	108.67
3	B	1304	SER	N-CA-C	9.06	122.20	109.15
3	A	1305	LEU	N-CA-C	-8.49	98.28	110.59
3	B	1305	LEU	N-CA-C	-7.12	100.26	110.59
3	A	1333	VAL	N-CA-C	6.11	116.28	110.53
3	B	1241	ILE	N-CA-C	-6.10	104.56	110.42
3	A	1390	CYS	N-CA-C	6.04	117.54	111.07
3	B	1271	ILE	N-CA-C	5.96	116.89	108.36
3	B	1333	VAL	N-CA-C	5.89	116.08	110.42
3	A	1135	ASP	N-CA-C	-5.74	106.44	113.50
3	B	1390	CYS	N-CA-C	5.68	117.15	111.07
3	B	1323	LYS	N-CA-C	5.39	117.58	111.11
3	A	1076	ASP	N-CA-C	-5.36	106.87	113.41
3	A	1053	LYS	N-CA-C	5.28	118.51	111.75
3	B	1384	HIS	N-CA-C	5.24	118.13	111.69
3	A	1385	GLN	N-CA-C	5.24	118.73	111.71
3	B	1392	LYS	N-CA-C	-5.21	105.68	111.36
3	A	1384	HIS	N-CA-C	5.15	116.97	111.36
3	A	1323	LYS	N-CA-C	5.07	117.21	111.02

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	268	0	149	9	0
2	D	181	0	104	7	0
3	A	3218	0	3158	79	0
3	B	3069	0	3020	104	0
4	A	14	0	0	0	0
4	B	14	0	0	0	0
4	D	7	0	0	0	0
5	A	25	0	11	0	0
5	B	25	0	11	1	0
6	A	18	0	9	0	0
6	B	6	0	3	0	0
7	B	1	0	0	1	0
8	A	266	0	0	9	0
8	B	199	0	0	8	0
8	C	21	0	0	0	0
8	D	8	0	0	0	0
All	All	7340	0	6465	193	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 14.

All (193) 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:B:1002:ARG:HH21	3:B:1002:ARG:HB3	1.13	1.11
3:A:1306:GLU:H	3:A:1309:HIS:HD2	1.07	0.97
3:B:1306:GLU:H	3:B:1309:HIS:HD2	1.04	0.95
3:B:1002:ARG:HB3	3:B:1002:ARG:NH2	1.85	0.92
3:A:1113:GLN:HE21	3:A:1115:ASP:H	1.16	0.87
3:A:1237:SER:HB2	3:A:1238:PRO:HD2	1.62	0.82
3:A:1306:GLU:H	3:A:1309:HIS:CD2	1.96	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:1306:GLU:H	3:B:1309:HIS:CD2	1.96	0.79
3:B:1113:GLN:HE21	3:B:1115:ASP:H	1.30	0.79
3:B:1002:ARG:HH21	3:B:1002:ARG:CB	1.95	0.78
3:B:1175:GLN:O	3:B:1177:PRO:HD3	1.84	0.77
3:B:1357:PHE:CE2	3:B:1361:LYS:HD2	2.19	0.77
3:A:1323:LYS:HE2	3:A:1327:GLU:OE1	1.87	0.74
3:B:1067:LYS:HE3	3:B:1067:LYS:HA	1.68	0.73
3:A:1332:ARG:H	3:A:1384:HIS:HD2	1.37	0.72
1:C:8:DA:H8	1:C:8:DA:H5'	1.56	0.70
3:B:1332:ARG:H	3:B:1384:HIS:HD2	1.44	0.66
1:C:3:DT:C3'	1:C:4:DA:H5'	2.26	0.65
3:A:1040:LYS:HB2	3:A:1040:LYS:NZ	2.12	0.65
3:B:1009:ARG:HH12	3:B:1044:PHE:HE2	1.43	0.65
3:B:1336:THR:OG1	3:B:1341:HIS:HE1	1.79	0.65
3:A:1354:GLU:O	3:A:1358:GLU:HG2	1.98	0.64
3:A:1242:ALA:O	3:A:1246:LYS:HG2	1.98	0.63
3:A:1313:GLY:HA3	3:A:1381:LEU:HD12	1.81	0.63
2:D:1:DC:H2'	2:D:2:DT:H72	1.80	0.62
3:A:1184:ARG:HG3	3:A:1188:TRP:HB2	1.81	0.62
3:A:1099:LEU:O	3:A:1103:ILE:HG13	1.98	0.62
3:B:1241:ILE:O	3:B:1245:GLU:HG3	2.00	0.61
3:A:1044:PHE:O	3:A:1047:THR:HG23	2.01	0.60
3:B:1286:LYS:NZ	3:B:1286:LYS:HB2	2.16	0.60
3:A:1322:TRP:CD1	3:A:1324:SER:HG	2.20	0.60
3:A:1113:GLN:NE2	3:A:1115:ASP:H	1.94	0.59
1:C:8:DA:H5'	1:C:8:DA:C8	2.38	0.59
3:B:1297:ASN:HD22	3:B:1298:GLN:H	1.51	0.59
3:B:1359:ARG:HH12	3:B:1363:LEU:HD21	1.68	0.59
3:B:1313:GLY:HA3	3:B:1381:LEU:HD12	1.83	0.59
3:B:1122:ARG:HG2	3:B:1122:ARG:HH21	1.68	0.58
3:A:1196:MET:HE3	3:A:1367:ARG:NH1	2.18	0.58
3:A:1218:HIS:HA	3:A:1222:LEU:HB2	1.85	0.58
2:D:5:DC:H2''	2:D:6:DT:H71	1.86	0.57
3:A:1357:PHE:O	3:A:1361:LYS:HG3	2.05	0.56
1:C:12:DA:H4'	3:B:1123:ARG:HE	1.70	0.56
3:B:1177:PRO:HB3	3:B:1390:CYS:HB2	1.88	0.56
3:A:1052:HIS:HD2	3:A:1056:SER:OG	1.88	0.56
3:B:1297:ASN:HD22	3:B:1298:GLN:N	2.04	0.55
3:A:1373:GLU:HA	3:A:1376:LYS:HD3	1.88	0.55
3:B:1223:LYS:HB3	3:B:1224:PRO:HD3	1.88	0.55
3:B:1050:HIS:HD2	8:B:95:HOH:O	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:1322:TRP:CD1	3:B:1324:SER:HG	2.24	0.55
3:B:1206:THR:HG22	3:B:1209:LYS:HG2	1.87	0.55
3:B:1023:GLN:HE22	3:B:1081:ASN:HD21	1.56	0.54
3:A:1191:VAL:HG23	8:A:74:HOH:O	2.06	0.54
3:A:1262:ASN:H	3:A:1270:GLN:NE2	2.05	0.54
3:B:1177:PRO:HB2	3:B:1387:SER:HA	1.89	0.54
3:B:1301:LEU:HD22	3:B:1329:LEU:HD21	1.89	0.54
3:B:1296:LEU:HD22	3:B:1296:LEU:N	2.23	0.54
3:B:1359:ARG:NH1	3:B:1363:LEU:HD21	2.22	0.54
3:A:1121:LEU:HB3	3:A:1150:LEU:HD11	1.89	0.53
3:A:1332:ARG:H	3:A:1384:HIS:CD2	2.23	0.53
3:B:1312:LEU:HD12	3:B:1321:PHE:HE1	1.73	0.53
3:B:1206:THR:CG2	3:B:1209:LYS:HG2	2.38	0.53
3:B:1336:THR:OG1	3:B:1341:HIS:CE1	2.61	0.53
3:B:1219:GLU:HG3	8:B:312:HOH:O	2.08	0.53
3:A:1223:LYS:HB3	3:A:1224:PRO:HD3	1.91	0.53
3:B:1119:LEU:HD12	3:B:1119:LEU:H	1.74	0.53
3:B:1332:ARG:H	3:B:1384:HIS:CD2	2.25	0.52
3:B:1097:LYS:HG2	3:B:1129:GLU:HG3	1.91	0.52
3:A:1336:THR:OG1	3:A:1341:HIS:CE1	2.63	0.52
3:B:1128:GLU:HB3	3:B:1129:GLU:OE1	2.10	0.52
3:B:1184:ARG:HG3	3:B:1188:TRP:HB2	1.92	0.52
3:B:1331:PHE:CE2	3:B:1333:VAL:HB	2.45	0.52
3:B:1009:ARG:NH1	3:B:1044:PHE:CE2	2.78	0.52
3:A:1346:ILE:HD11	3:A:1363:LEU:HD12	1.91	0.52
3:B:1295:LYS:HB2	8:B:468:HOH:O	2.09	0.52
3:B:1296:LEU:HB2	3:B:1301:LEU:HD21	1.92	0.52
3:B:1089:GLN:HG2	8:B:85:HOH:O	2.10	0.52
2:D:1:DC:H5"	3:B:1245:GLU:OE1	2.09	0.51
3:A:1260:LYS:NZ	3:A:1260:LYS:HB3	2.24	0.51
3:B:1040:LYS:HG3	3:B:1060:ILE:HG12	1.92	0.51
3:A:1372:ARG:O	3:A:1376:LYS:HG3	2.10	0.51
3:B:1286:LYS:HB2	3:B:1286:LYS:HZ2	1.75	0.51
3:B:1047:THR:O	3:B:1052:HIS:HE1	1.93	0.51
3:B:1012:GLU:HG2	3:B:1044:PHE:CE2	2.46	0.51
3:B:1119:LEU:HD12	3:B:1119:LEU:N	2.25	0.51
3:B:1191:VAL:HG23	8:B:99:HOH:O	2.10	0.51
3:A:1148:LYS:O	3:A:1152:LYS:HD2	2.11	0.51
3:B:1179:ASP:OD1	3:B:1181:VAL:HG22	2.11	0.51
3:B:1122:ARG:HG2	3:B:1122:ARG:NH2	2.26	0.50
3:B:1257:GLU:HG2	3:B:1257:GLU:O	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:1023:GLN:HE22	3:A:1081:ASN:HD21	1.58	0.50
3:A:1357:PHE:CE2	3:A:1361:LYS:HD2	2.47	0.50
3:A:1010:GLY:HA2	3:A:1039:ALA:O	2.11	0.50
3:B:1189:LYS:HD3	8:B:289:HOH:O	2.10	0.50
3:A:1336:THR:OG1	3:A:1341:HIS:HE1	1.94	0.50
3:B:1046:ARG:NH2	3:B:1235:GLU:OE2	2.45	0.50
3:A:1198:ILE:HD12	8:A:341:HOH:O	2.11	0.49
3:B:1140:HIS:CE1	3:B:1176:PRO:HD3	2.46	0.49
3:A:1128:GLU:O	3:A:1132:ARG:HG3	2.12	0.49
3:A:1254:GLY:H	3:A:1257:GLU:CD	2.20	0.49
3:B:1273:ASP:OD2	3:B:1274:CME:N	2.45	0.49
3:A:1040:LYS:HB2	3:A:1040:LYS:HZ1	1.76	0.49
3:A:1313:GLY:HA3	3:A:1381:LEU:CD1	2.43	0.49
3:B:1260:LYS:HB3	3:B:1260:LYS:NZ	2.27	0.49
3:B:1254:GLY:C	3:B:1256:ARG:H	2.19	0.49
3:A:1309:HIS:HE1	8:A:8:HOH:O	1.95	0.49
3:B:1213:GLN:HG2	3:B:1322:TRP:CZ3	2.47	0.49
3:A:1053:LYS:HE2	8:A:344:HOH:O	2.12	0.49
3:B:1052:HIS:HD2	3:B:1056:SER:OG	1.96	0.49
3:B:1002:ARG:HD3	8:B:155:HOH:O	2.14	0.48
3:A:1050:HIS:HD2	8:A:62:HOH:O	1.97	0.48
3:B:1177:PRO:HG2	3:B:1391:PHE:CE1	2.48	0.48
3:B:1208:TRP:HA	3:B:1295:LYS:HE3	1.95	0.48
3:A:1236:ARG:HG3	3:A:1252:TYR:CZ	2.49	0.47
3:B:1199:ASN:HB3	3:B:1291:TYR:CE2	2.49	0.47
2:D:1:DC:C2'	2:D:2:DT:H72	2.43	0.47
3:A:1047:THR:O	3:A:1052:HIS:HE1	1.97	0.47
1:C:11:DT:O4'	3:B:1119:LEU:HD22	2.15	0.47
3:A:1199:ASN:HB3	3:A:1291:TYR:CE2	2.50	0.47
3:B:1012:GLU:HB2	7:B:4001:CL:CL	2.51	0.47
3:B:1113:GLN:NE2	3:B:1115:ASP:H	2.04	0.47
3:A:1231:MET:HB2	3:A:1271:ILE:CD1	2.45	0.47
3:B:1296:LEU:HB2	3:B:1301:LEU:CG	2.45	0.47
3:A:1326:GLY:HA3	3:A:1347:TRP:CZ3	2.50	0.47
3:A:1113:GLN:HE21	3:A:1115:ASP:N	1.99	0.46
3:B:1067:LYS:HA	3:B:1067:LYS:CE	2.41	0.46
3:B:1251:GLU:HG2	3:B:1253:TYR:OH	2.14	0.46
3:B:1303:ARG:O	3:B:1384:HIS:CE1	2.67	0.46
3:A:1180:ILE:HG23	3:A:1314:ALA:HB2	1.96	0.46
1:C:3:DT:H3'	1:C:4:DA:H5'	1.95	0.46
1:C:3:DT:H3'	1:C:4:DA:C5'	2.44	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:1140:HIS:HE1	3:B:1176:PRO:HD3	1.79	0.46
3:A:1299:LYS:HE2	8:A:254:HOH:O	2.17	0.45
3:B:1372:ARG:O	3:B:1376:LYS:HG3	2.16	0.45
3:B:1039:ALA:HA	3:B:1059:VAL:O	2.16	0.45
3:B:1086:THR:HA	3:B:1125:LEU:CD1	2.47	0.45
3:A:1199:ASN:ND2	3:A:1291:TYR:OH	2.49	0.44
3:B:1297:ASN:ND2	3:B:1298:GLN:N	2.65	0.44
1:C:3:DT:H5'	3:A:1207:THR:HG21	1.98	0.44
3:A:1086:THR:HA	3:A:1125:LEU:HG	1.99	0.44
3:A:1129:GLU:H	3:A:1129:GLU:CD	2.23	0.44
3:B:1129:GLU:H	3:B:1129:GLU:CD	2.25	0.44
3:A:1065:TYR:CZ	3:A:1098:LYS:HD2	2.52	0.44
2:D:8:DA:C2	3:A:1238:PRO:HB2	2.53	0.44
3:B:1085:ALA:O	3:B:1088:VAL:HG22	2.18	0.44
3:B:1086:THR:HA	3:B:1125:LEU:HD11	2.00	0.43
3:B:1113:GLN:NE2	3:B:1115:ASP:HB2	2.33	0.43
3:B:1369:LEU:CD1	3:B:1372:ARG:HH11	2.31	0.43
3:B:1254:GLY:C	3:B:1256:ARG:N	2.75	0.43
3:A:1040:LYS:NZ	3:A:1040:LYS:CB	2.80	0.43
3:A:1306:GLU:N	3:A:1309:HIS:HD2	1.92	0.43
3:A:1038:TYR:CZ	3:A:1058:PRO:HB3	2.54	0.43
3:B:1206:THR:HG22	3:B:1209:LYS:CG	2.49	0.43
3:A:1002:ARG:NH2	8:A:87:HOH:O	2.51	0.43
3:B:1262:ASN:H	3:B:1270:GLN:NE2	2.16	0.43
3:B:1275:TYR:CE1	5:B:2001:UDP:C2	3.06	0.43
3:B:1301:LEU:HD22	3:B:1329:LEU:CG	2.49	0.43
3:A:1231:MET:HB2	3:A:1271:ILE:HD12	1.99	0.43
1:C:3:DT:C3'	1:C:4:DA:C5'	2.97	0.42
3:A:1254:GLY:H	3:A:1257:GLU:CG	2.32	0.42
3:A:1258:ILE:O	3:A:1261:MET:HB2	2.19	0.42
3:A:1244:LYS:HA	8:A:399:HOH:O	2.19	0.42
3:B:1329:LEU:O	3:B:1337:PRO:HA	2.19	0.42
3:A:1262:ASN:C	3:A:1262:ASN:HD22	2.28	0.42
3:B:1331:PHE:HE1	3:B:1384:HIS:HB2	1.85	0.42
2:D:2:DT:OP2	3:A:1043:SER:CB	2.68	0.42
2:D:5:DC:C2'	2:D:6:DT:H71	2.48	0.42
3:B:1301:LEU:HD12	3:B:1328:ASN:HB2	2.02	0.42
3:B:1138:PHE:HA	3:B:1172:TYR:O	2.20	0.41
3:B:1264:ALA:HB1	3:B:1265:PRO:HD2	2.02	0.41
3:A:1208:TRP:O	3:A:1294:SER:HA	2.20	0.41
3:B:1309:HIS:HE1	8:B:51:HOH:O	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:1331:PHE:CE1	3:B:1384:HIS:HB2	2.55	0.41
3:A:1359:ARG:NH1	3:A:1363:LEU:HD21	2.36	0.41
3:B:1036:LEU:C	3:B:1036:LEU:HD13	2.46	0.41
3:A:1138:PHE:HA	3:A:1172:TYR:O	2.20	0.41
3:B:1038:TYR:CZ	3:B:1058:PRO:HB3	2.55	0.41
3:B:1373:GLU:O	3:B:1376:LYS:HB2	2.21	0.41
3:A:1039:ALA:HA	3:A:1059:VAL:O	2.21	0.41
3:A:1306:GLU:O	3:A:1309:HIS:HB2	2.21	0.41
3:B:1390:CYS:O	3:B:1394:GLN:HG3	2.21	0.41
3:A:1008:ALA:HB3	3:A:1083:VAL:HA	2.03	0.41
3:A:1017:THR:O	3:A:1021:LEU:HG	2.21	0.41
3:B:1178:MET:HE3	3:B:1183:VAL:HG21	2.03	0.41
3:A:1359:ARG:HH11	3:A:1363:LEU:HG	1.86	0.41
3:A:1363:LEU:HD21	3:A:1373:GLU:HG3	2.03	0.41
3:B:1113:GLN:HE21	3:B:1115:ASP:N	2.07	0.41
3:A:1253:TYR:OH	3:A:1270:GLN:NE2	2.53	0.40
3:A:1066:ASP:HB3	8:A:77:HOH:O	2.20	0.40
3:A:1176:PRO:HA	3:A:1177:PRO:HD3	1.91	0.40
3:B:1216:ASP:OD1	3:B:1353:MET:HE1	2.22	0.40
3:A:1248:ILE:HA	3:A:1249:PRO:HD3	1.89	0.40
3:B:1218:HIS:HE1	3:B:1267:GLN:O	2.04	0.40
3:B:1262:ASN:C	3:B:1262:ASN:HD22	2.28	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	
3	A	385/402 (96%)	365 (95%)	20 (5%)	0	100	100
3	B	367/402 (91%)	354 (96%)	13 (4%)	0	100	100
All	All	752/804 (94%)	719 (96%)	33 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	354/365 (97%)	352 (99%)	2 (1%)	78	81
3	B	338/365 (93%)	334 (99%)	4 (1%)	63	63
All	All	692/730 (95%)	686 (99%)	6 (1%)	70	73

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

Mol	Chain	Res	Type
3	A	1040	LYS
3	A	1129	GLU
3	B	1002	ARG
3	B	1061	LEU
3	B	1067	LYS
3	B	1129	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (46) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	1023	GLN
3	A	1050	HIS
3	A	1052	HIS
3	A	1094	ASN
3	A	1113	GLN
3	A	1143	ASN
3	A	1175	GLN
3	A	1199	ASN
3	A	1218	HIS
3	A	1266	ASN
3	A	1267	GLN
3	A	1270	GLN
3	A	1277	ASN

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Mol	Chain	Res	Type
3	A	1309	HIS
3	A	1328	ASN
3	A	1341	HIS
3	A	1351	ASN
3	A	1383	GLN
3	A	1384	HIS
3	A	1394	GLN
3	B	1023	GLN
3	B	1030	ASN
3	B	1050	HIS
3	B	1052	HIS
3	B	1089	GLN
3	B	1094	ASN
3	B	1102	ASN
3	B	1113	GLN
3	B	1114	HIS
3	B	1124	ASN
3	B	1140	HIS
3	B	1175	GLN
3	B	1199	ASN
3	B	1218	HIS
3	B	1262	ASN
3	B	1266	ASN
3	B	1267	GLN
3	B	1270	GLN
3	B	1297	ASN
3	B	1302	GLN
3	B	1309	HIS
3	B	1328	ASN
3	B	1341	HIS
3	B	1351	ASN
3	B	1384	HIS
3	B	1394	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

5 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	$\# Z > 2$
3	CME	B	1014	3	8,9,10	1.08	1 (12%)	6,9,11	0.38	0
3	CME	B	1274	3	8,9,10	1.07	1 (12%)	6,9,11	0.48	0
3	CME	A	1274	3	8,9,10	1.08	1 (12%)	6,9,11	0.42	0
1	5HU	C	7	1	19,22,23	0.38	0	25,31,34	0.77	1 (4%)
3	CME	A	1014	3	8,9,10	1.11	1 (12%)	6,9,11	0.57	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	CME	B	1014	3	-	1/5/8/10	-
3	CME	B	1274	3	-	1/5/8/10	-
3	CME	A	1274	3	-	1/5/8/10	-
1	5HU	C	7	1	-	0/9/23/24	0/2/2/2
3	CME	A	1014	3	-	2/5/8/10	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1014	CME	OH-CZ	-2.94	1.27	1.42
3	B	1014	CME	OH-CZ	-2.88	1.27	1.42
3	B	1274	CME	OH-CZ	-2.78	1.27	1.42
3	A	1274	CME	OH-CZ	-2.76	1.28	1.42

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	7	5HU	C5A-C5-C6	2.90	123.37	119.25

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1014	CME	SD-CE-CZ-OH
3	B	1014	CME	CE-SD-SG-CB
3	A	1014	CME	CZ-CE-SD-SG
3	B	1274	CME	CZ-CE-SD-SG
3	A	1274	CME	CA-CB-SG-SD

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1274	CME	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 1 is monoatomic - leaving 11 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
6	GOL	B	5004	-	5,5,5	4.50	5 (100%)	5,5,5	2.23	2 (40%)
4	NCO	D	3003	-	6,6,6	1.49	0	-		
4	NCO	B	3004	-	6,6,6	1.47	0	-		
6	GOL	A	5003	-	5,5,5	4.49	5 (100%)	5,5,5	2.23	2 (40%)
4	NCO	B	3002	-	6,6,6	1.52	0	-		
4	NCO	A	3005	-	6,6,6	1.47	0	-		
6	GOL	A	5002	-	5,5,5	4.50	5 (100%)	5,5,5	2.19	2 (40%)
5	UDP	A	3006	-	25,26,26	0.96	2 (8%)	38,40,40	0.60	1 (2%)
4	NCO	A	3001	-	6,6,6	1.51	0	-		
6	GOL	A	5001	-	5,5,5	4.49	5 (100%)	5,5,5	2.24	2 (40%)
5	UDP	B	2001	-	25,26,26	0.96	1 (4%)	38,40,40	0.59	1 (2%)

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
6	GOL	B	5004	-	-	2/4/4/4	-
6	GOL	A	5003	-	-	2/4/4/4	-
6	GOL	A	5002	-	-	2/4/4/4	-
5	UDP	A	3006	-	-	4/16/32/32	0/2/2/2
6	GOL	A	5001	-	-	3/4/4/4	-
5	UDP	B	2001	-	-	4/16/32/32	0/2/2/2

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	5002	GOL	C1-C2	-5.59	1.30	1.51
6	B	5004	GOL	C3-C2	-5.57	1.30	1.51
6	A	5002	GOL	C3-C2	-5.56	1.30	1.51
6	A	5001	GOL	C1-C2	-5.54	1.30	1.51
6	A	5003	GOL	C3-C2	-5.54	1.30	1.51
6	A	5001	GOL	C3-C2	-5.53	1.30	1.51
6	B	5004	GOL	C1-C2	-5.50	1.30	1.51
6	A	5003	GOL	C1-C2	-5.48	1.30	1.51
6	B	5004	GOL	O1-C1	4.01	1.59	1.42
6	A	5003	GOL	O1-C1	4.00	1.59	1.42
6	A	5001	GOL	O3-C3	3.99	1.59	1.42
6	A	5001	GOL	O1-C1	3.95	1.59	1.42
6	B	5004	GOL	O3-C3	3.95	1.59	1.42
6	A	5002	GOL	O1-C1	3.92	1.58	1.42
6	A	5002	GOL	O3-C3	3.91	1.58	1.42
6	A	5003	GOL	O3-C3	3.91	1.58	1.42
5	B	2001	UDP	PA-O3A	3.82	1.63	1.59
5	A	3006	UDP	PA-O3A	3.64	1.63	1.59
6	A	5003	GOL	O2-C2	-2.96	1.34	1.43
6	B	5004	GOL	O2-C2	-2.87	1.35	1.43
6	A	5002	GOL	O2-C2	-2.85	1.35	1.43
6	A	5001	GOL	O2-C2	-2.82	1.35	1.43
5	A	3006	UDP	PB-O2B	-2.02	1.47	1.54

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	5003	GOL	O1-C1-C2	3.51	126.20	110.38
6	B	5004	GOL	O1-C1-C2	3.49	126.11	110.38
6	A	5001	GOL	O3-C3-C2	3.48	126.03	110.38
6	A	5001	GOL	O1-C1-C2	3.45	125.93	110.38
6	A	5003	GOL	O3-C3-C2	3.42	125.77	110.38
6	B	5004	GOL	O3-C3-C2	3.41	125.73	110.38
6	A	5002	GOL	O3-C3-C2	3.41	125.73	110.38
6	A	5002	GOL	O1-C1-C2	3.38	125.61	110.38
5	A	3006	UDP	O2B-PB-O1B	2.21	119.44	110.83
5	B	2001	UDP	O2B-PB-O1B	2.17	119.27	110.83

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	3006	UDP	PA-O3A-PB-O3B
5	B	2001	UDP	PA-O3A-PB-O3B
6	A	5001	GOL	C1-C2-C3-O3
6	B	5004	GOL	O1-C1-C2-C3
6	A	5001	GOL	O2-C2-C3-O3
6	A	5002	GOL	C1-C2-C3-O3
6	A	5003	GOL	O1-C1-C2-C3
6	A	5002	GOL	O2-C2-C3-O3
6	A	5003	GOL	O1-C1-C2-O2
6	B	5004	GOL	O1-C1-C2-O2
5	B	2001	UDP	PB-O3A-PA-O5'
6	A	5001	GOL	O1-C1-C2-O2
5	A	3006	UDP	C5'-O5'-PA-O3A
5	A	3006	UDP	PA-O3A-PB-O2B
5	A	3006	UDP	O4'-C4'-C5'-O5'
5	B	2001	UDP	O4'-C4'-C5'-O5'
5	B	2001	UDP	PA-O3A-PB-O1B

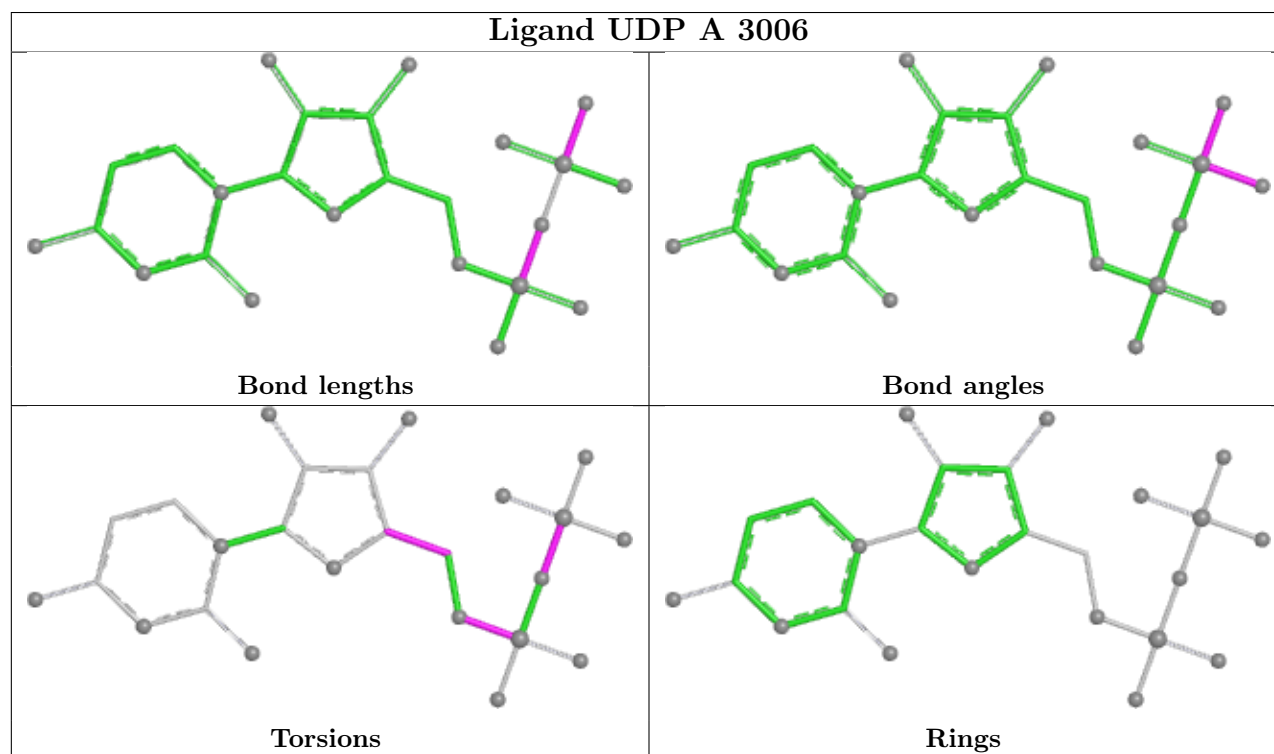
There are no ring outliers.

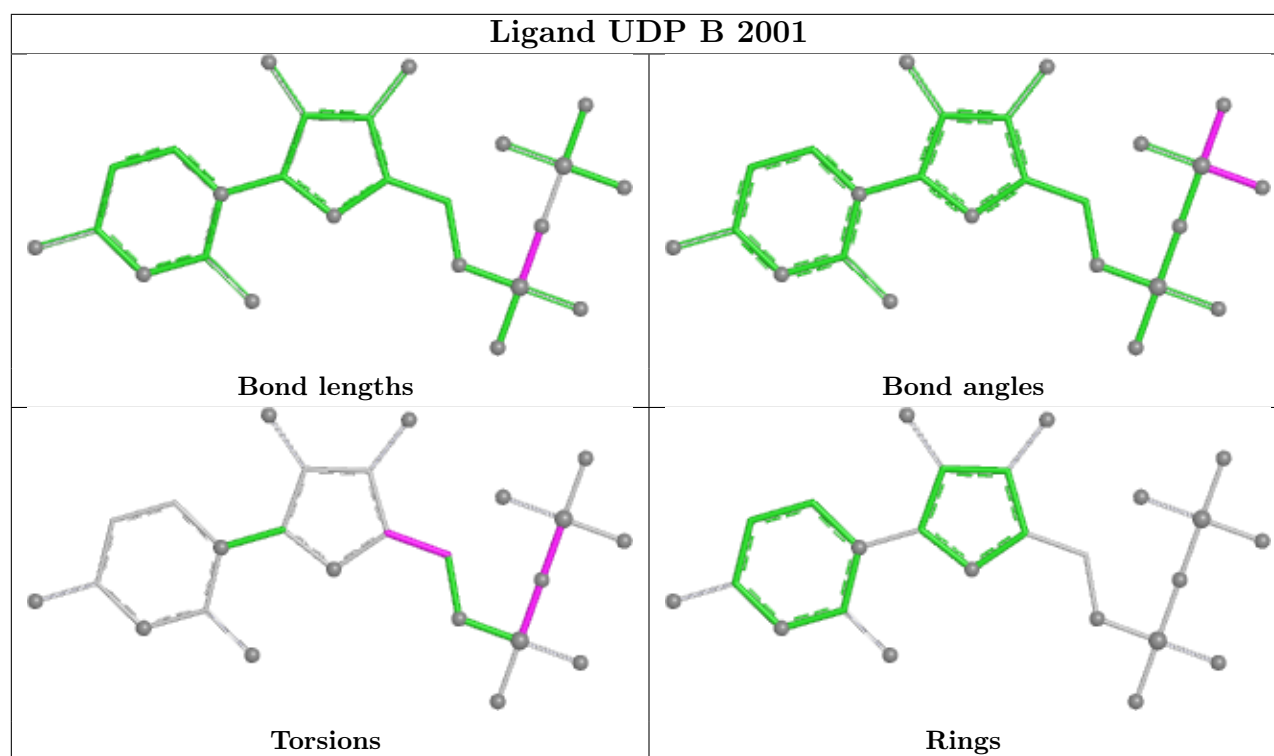
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	2001	UDP	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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	C	12/13 (92%)	0.56	1 (8%) 17 18	24, 46, 59, 74	0
2	D	9/9 (100%)	0.68	0 100 100	41, 47, 50, 50	0
3	A	389/402 (96%)	0.19	13 (3%) 49 52	11, 25, 43, 60	0
3	B	371/402 (92%)	0.40	16 (4%) 40 42	16, 30, 50, 67	0
All	All	781/826 (94%)	0.30	30 (3%) 44 47	11, 28, 48, 74	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	B	1301	LEU	4.0
3	B	1399	THR	3.3
3	B	1300	TYR	3.2
3	A	1258	ILE	3.1
3	B	1297	ASN	3.0
3	A	1000	SER	3.0
3	A	1252	TYR	2.9
3	B	1298	GLN	2.6
3	B	1258	ILE	2.5
3	B	1296	LEU	2.5
3	B	1256	ARG	2.4
3	B	1181	VAL	2.4
3	B	1132	ARG	2.3
3	A	1089	GLN	2.3
3	A	999	GLY	2.3
3	A	1251	GLU	2.3
1	C	1	DG	2.3
3	A	1266	ASN	2.3
3	B	1120	SER	2.2
3	B	1119	LEU	2.2
3	B	1171	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
3	B	1086	THR	2.2
3	A	1191	VAL	2.1
3	A	1365	SER	2.1
3	B	1340	SER	2.1
3	A	1262	ASN	2.1
3	A	1066	ASP	2.1
3	B	1400	LYS	2.0
3	A	1358	GLU	2.0
3	A	1255	ASN	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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	B-factors(Å ²)	Q<0.9
3	CME	A	1274	10/11	0.85	0.13	28,37,49,51	0
3	CME	B	1274	10/11	0.88	0.12	25,32,45,46	0
1	5HU	C	7	21/22	0.89	0.11	33,47,49,50	0
3	CME	B	1014	10/11	0.92	0.10	23,29,42,46	0
3	CME	A	1014	10/11	0.92	0.10	13,22,38,39	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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. 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	B-factors(Å ²)	Q<0.9
6	GOL	B	5004	6/6	0.58	0.16	67,68,68,69	0
6	GOL	A	5002	6/6	0.70	0.15	63,63,63,64	0
6	GOL	A	5001	6/6	0.78	0.14	72,72,72,72	0
4	NCO	B	3004	7/7	0.79	0.22	53,53,53,54	7

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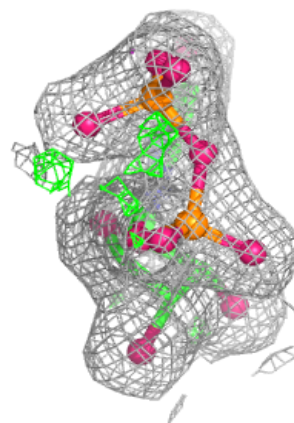
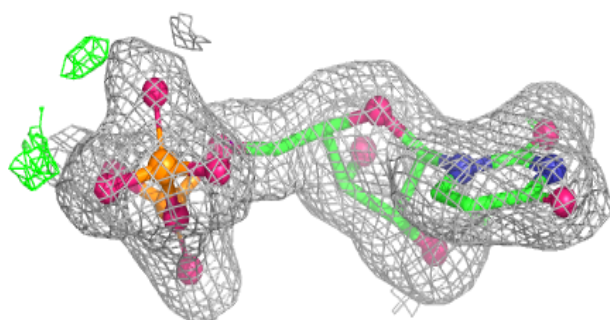
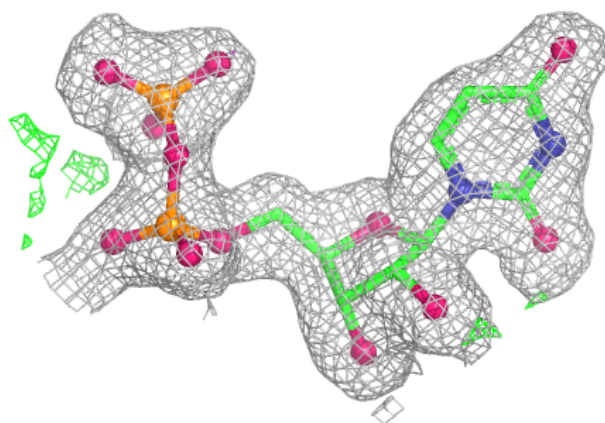
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	GOL	A	5003	6/6	0.80	0.14	62,63,63,64	0
4	NCO	A	3005	7/7	0.89	0.14	45,45,45,46	7
4	NCO	D	3003	7/7	0.93	0.11	58,58,59,59	0
4	NCO	B	3002	7/7	0.94	0.11	43,44,44,45	0
5	UDP	A	3006	25/25	0.97	0.06	13,16,20,22	0
5	UDP	B	2001	25/25	0.97	0.05	17,20,23,25	0
4	NCO	A	3001	7/7	0.97	0.08	43,44,44,45	0
7	CL	B	4001	1/1	0.98	0.06	36,36,36,36	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

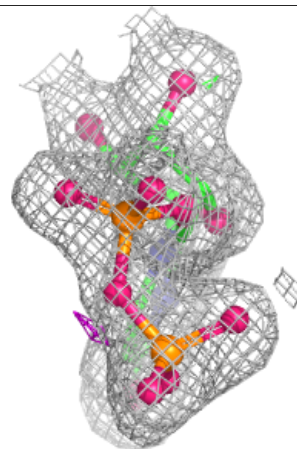
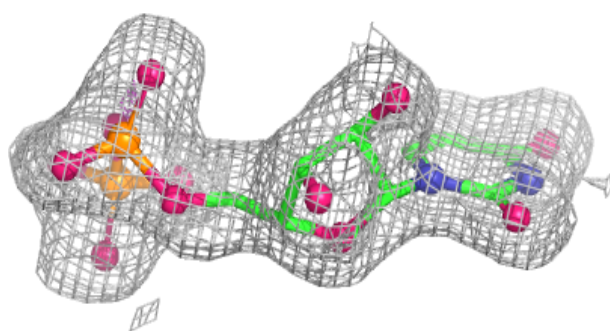
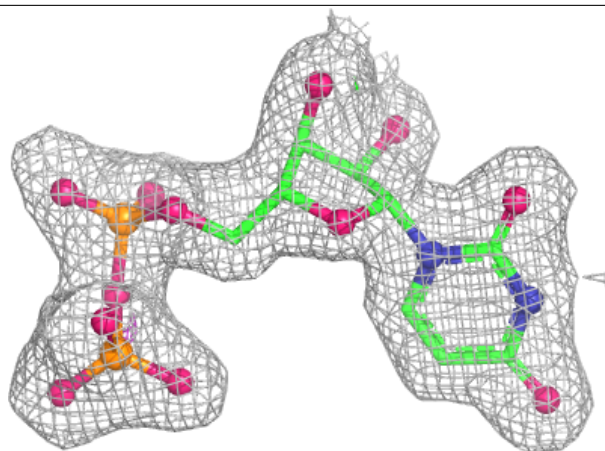
Electron density around UDP A 3006:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around UDP B 2001:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.