



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 8, 2023 – 06:04 PM JST

PDB ID : 8J30
Title : Crystal structure of ApNGT with Q469A and M218A mutations in complex with UDP-GLC
Authors : Feng, Y.; Hao, Z.; Guo, Q.; Zheng, J.; Da, L.; Peng, W.
Deposited on : 2023-04-15
Resolution : 2.89 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.35
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35

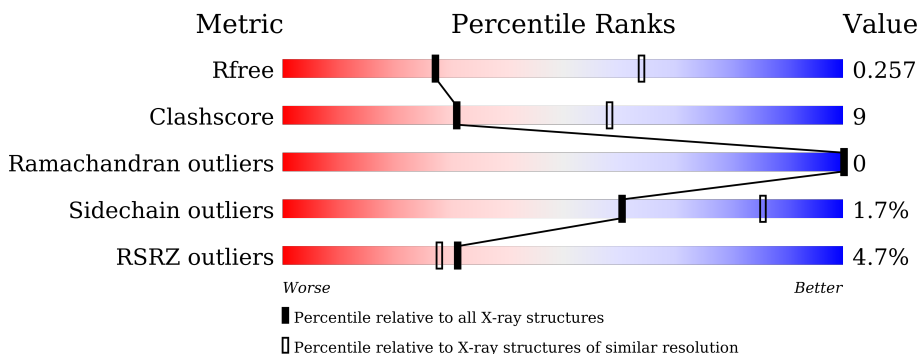
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

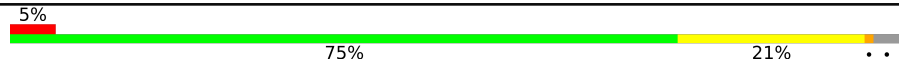

The reported resolution of this entry is 2.89 Å.

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}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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	A	628	 5% 75% 21% ..
1	B	628	 4% 76% 18% 5%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 9688 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.

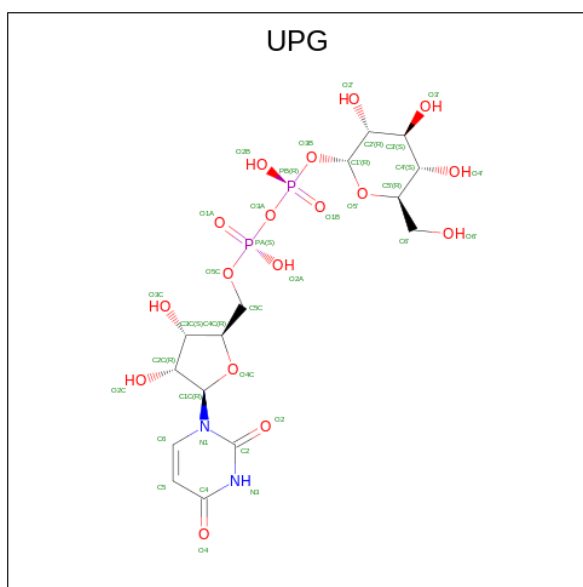
- Molecule 1 is a protein called UDP-glucose:protein N-beta-glucosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	608	4860	3106	832	898	24	0	0	0
1	B	596	4767	3050	816	878	23	0	0	0

There are 20 discrepancies between the modelled and reference sequences:

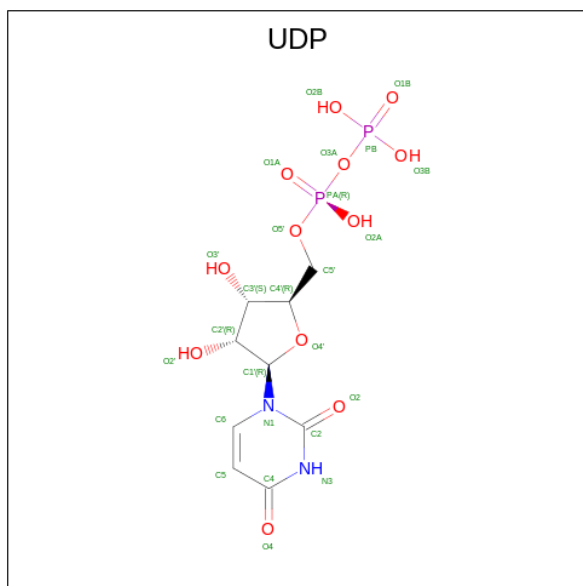
Chain	Residue	Modelled	Actual	Comment	Reference
A	218	ALA	MET	engineered mutation	UNP A3N2T3
A	469	ALA	GLN	engineered mutation	UNP A3N2T3
A	621	LEU	-	expression tag	UNP A3N2T3
A	622	GLU	-	expression tag	UNP A3N2T3
A	623	HIS	-	expression tag	UNP A3N2T3
A	624	HIS	-	expression tag	UNP A3N2T3
A	625	HIS	-	expression tag	UNP A3N2T3
A	626	HIS	-	expression tag	UNP A3N2T3
A	627	HIS	-	expression tag	UNP A3N2T3
A	628	HIS	-	expression tag	UNP A3N2T3
B	218	ALA	MET	engineered mutation	UNP A3N2T3
B	469	ALA	GLN	engineered mutation	UNP A3N2T3
B	621	LEU	-	expression tag	UNP A3N2T3
B	622	GLU	-	expression tag	UNP A3N2T3
B	623	HIS	-	expression tag	UNP A3N2T3
B	624	HIS	-	expression tag	UNP A3N2T3
B	625	HIS	-	expression tag	UNP A3N2T3
B	626	HIS	-	expression tag	UNP A3N2T3
B	627	HIS	-	expression tag	UNP A3N2T3
B	628	HIS	-	expression tag	UNP A3N2T3

- Molecule 2 is URIDINE-5'-DIPHOSPHATE-GLUCOSE (three-letter code: UPG) (formula: C₁₅H₂₄N₂O₁₇P₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	N	O			P
2	A	1	36	15	2	17	2	0	0

- Molecule 3 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: C₉H₁₄N₂O₁₂P₂) (labeled as "Ligand of Interest" by depositor).

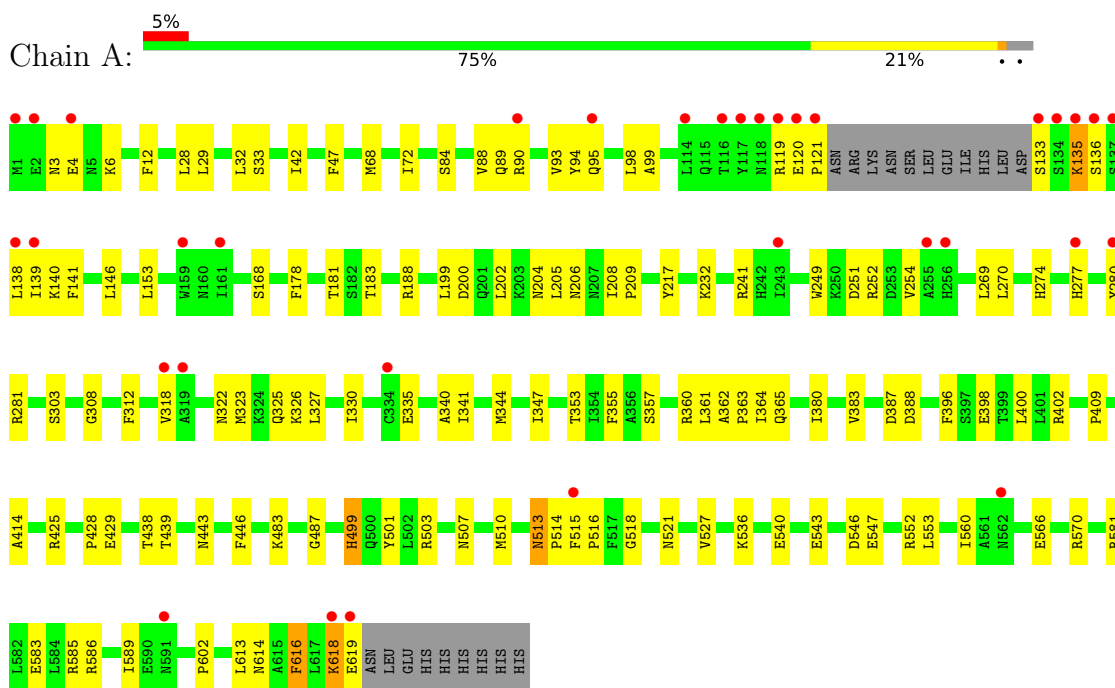


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	N	O			P
3	B	1	25	9	2	12	2	0	0

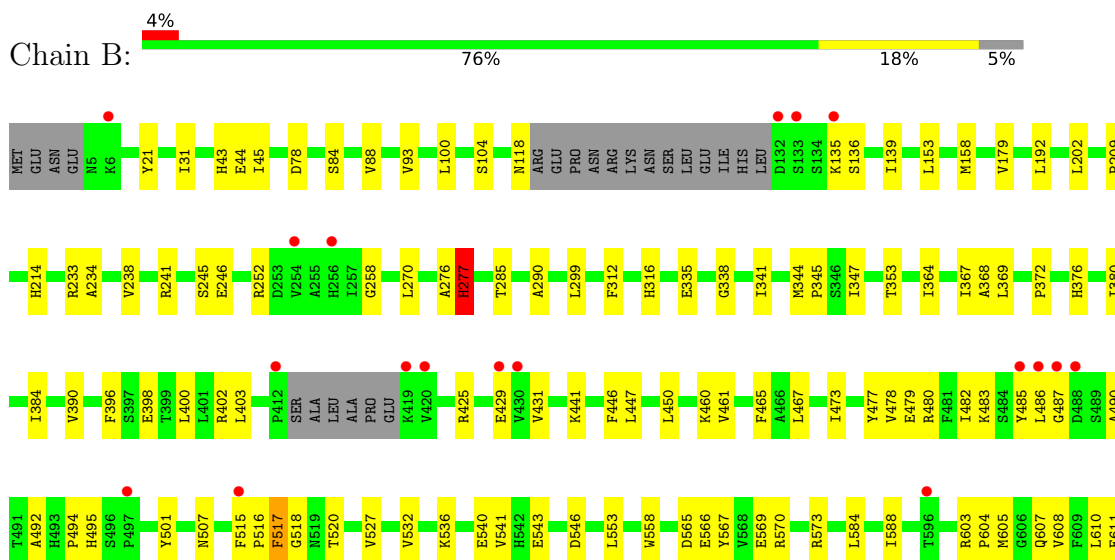
3 Residue-property plots i

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: UDP-glucose:protein N-beta-glucosyltransferase



- Molecule 1: UDP-glucose:protein N-beta-glucosyltransferase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	79.41Å 95.01Å 175.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.18 – 2.89 47.51 – 2.89	Depositor EDS
% Data completeness (in resolution range)	98.7 (47.18-2.89) 98.8 (47.51-2.89)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.64 (at 2.91Å)	Xtrriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.211 , 0.250 0.223 , 0.257	Depositor DCC
R_{free} test set	1528 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	51.8	Xtrriage
Anisotropy	0.055	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 43.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	9688	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.01% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: UDP, UPG

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	0.40	0/4980	0.59	0/6763
1	B	0.40	0/4884	0.61	1/6631 (0.0%)
All	All	0.40	0/9864	0.60	1/13394 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	277	HIS	N-CA-CB	5.20	119.96	110.60

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	4860	0	4772	90	2
1	B	4767	0	4680	84	2
2	A	36	0	22	3	0
3	B	25	0	11	1	0
All	All	9688	0	9485	174	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.

All (174) 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:178:PHE:HB2	1:A:439:THR:HB	1.37	1.05
1:B:483:LYS:HA	1:B:487:GLY:HA2	1.58	0.84
1:A:274:HIS:HB3	1:A:277:HIS:CD2	2.15	0.82
1:A:425:ARG:HD2	1:A:428:PRO:HB3	1.63	0.81
1:B:515:PHE:HD1	1:B:536:LYS:HD2	1.45	0.80
1:B:233:ARG:HG2	1:B:376:HIS:HB3	1.62	0.80
1:A:527:VAL:HG21	1:A:553:LEU:HD13	1.64	0.79
1:B:447:LEU:HB3	1:B:485:TYR:CG	2.23	0.73
1:B:258:GLY:N	1:B:338:GLY:HA3	2.05	0.72
1:A:585:ARG:O	1:A:589:ILE:HG13	1.90	0.71
1:B:483:LYS:HA	1:B:487:GLY:CA	2.23	0.69
1:A:483:LYS:HG2	1:A:487:GLY:HA2	1.75	0.69
1:A:90:ARG:O	1:A:93:VAL:HG22	1.94	0.67
1:A:280:TYR:HE2	1:A:308:GLY:HA2	1.57	0.67
1:A:323:MET:HG2	1:A:355:PHE:CD2	2.30	0.67
1:B:447:LEU:HD12	1:B:485:TYR:HB2	1.77	0.66
1:A:200:ASP:O	1:A:241:ARG:NH2	2.28	0.65
1:A:521:ASN:HD22	2:A:701:UPG:PB	2.18	0.65
1:A:614:ASN:O	1:A:618:LYS:HE3	1.99	0.62
1:A:68:MET:O	1:A:72:ILE:HG13	2.00	0.62
1:A:280:TYR:CE2	1:A:308:GLY:HA2	2.34	0.62
1:A:616:PHE:O	1:A:619:GLU:HG3	1.99	0.61
1:A:90:ARG:HE	1:A:94:TYR:HE2	1.48	0.61
1:B:565:ASP:O	1:B:569:GLU:HG3	2.01	0.61
1:A:341:ILE:HD13	1:A:364:ILE:HB	1.82	0.60
1:B:479:GLU:O	1:B:483:LYS:HG2	2.01	0.60
1:B:380:ILE:HG22	1:B:398:GLU:HG2	1.84	0.60
1:B:566:GLU:O	1:B:570:ARG:HG2	2.02	0.59
1:A:388:ASP:OD1	1:A:388:ASP:N	2.34	0.58
1:B:494:PRO:O	1:B:495:HIS:C	2.41	0.58
1:A:443:ASN:ND2	1:A:516:PRO:HG3	2.19	0.58
1:B:483:LYS:CA	1:B:487:GLY:HA2	2.32	0.58
1:A:501:TYR:CE2	2:A:701:UPG:H2C	2.38	0.58
1:A:84:SER:O	1:A:88:VAL:HG23	2.04	0.57
1:B:465:PHE:HE1	1:B:490:ALA:HB1	1.69	0.57
1:A:322:ASN:OD1	1:A:325:GLN:HG3	2.04	0.57
1:A:204:ASN:HD22	1:A:206:ASN:H	1.51	0.57
1:B:390:VAL:HG21	1:B:402:ARG:HG2	1.86	0.57
1:A:387:ASP:OD1	1:A:402:ARG:NH2	2.38	0.57
1:A:29:LEU:HD22	1:A:90:ARG:CD	2.35	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:318:VAL:HG12	1:A:326:LYS:HG2	1.86	0.56
1:B:441:LYS:HD3	1:B:517:PHE:O	2.06	0.56
1:A:547:GLU:HG3	1:A:560:ILE:HG21	1.87	0.56
1:A:33:SER:HB2	1:A:90:ARG:HH21	1.72	0.55
1:B:515:PHE:CD1	1:B:536:LYS:HD2	2.35	0.55
1:B:607:GLN:O	1:B:611:GLU:HG3	2.07	0.55
1:A:205:LEU:HD23	1:A:208:ILE:HD11	1.88	0.55
1:B:482:ILE:HD12	1:B:492:ALA:HB2	1.88	0.55
1:A:521:ASN:N	2:A:701:UPG:O2B	2.37	0.54
1:B:617:LEU:HB2	1:B:618:LYS:HD2	1.89	0.54
1:B:252:ARG:NE	1:B:335:GLU:OE1	2.40	0.54
1:B:135:LYS:HE3	1:B:139:ILE:HD11	1.90	0.54
1:A:360:ARG:HH11	1:A:365:GLN:HG3	1.72	0.54
1:B:290:ALA:HB2	1:B:605:MET:HE2	1.90	0.54
1:B:429:GLU:O	1:B:460:LYS:HB2	2.08	0.54
1:A:518:GLY:HA3	1:A:546:ASP:OD1	2.07	0.54
1:A:513:ASN:HB3	1:A:514:PRO:HD2	1.91	0.53
1:B:84:SER:O	1:B:88:VAL:HG23	2.09	0.53
1:B:613:LEU:O	1:B:617:LEU:HG	2.09	0.53
1:B:425:ARG:NH2	1:B:507:ASN:O	2.40	0.53
1:B:518:GLY:HA3	1:B:546:ASP:OD1	2.09	0.53
1:B:483:LYS:HA	1:B:487:GLY:H	1.74	0.52
1:A:515:PHE:CD2	1:A:516:PRO:HD3	2.44	0.52
1:B:43:HIS:O	1:B:44:GLU:HB2	2.09	0.52
1:B:532:VAL:HG21	1:B:584:LEU:HD23	1.91	0.52
1:B:515:PHE:HB2	1:B:516:PRO:HD3	1.92	0.52
1:A:119:ARG:HD2	1:A:133:SER:HB3	1.91	0.52
1:B:446:PHE:HA	1:B:515:PHE:CE2	2.44	0.52
1:B:483:LYS:HA	1:B:487:GLY:N	2.24	0.52
1:A:120:GLU:N	1:A:121:PRO:HD3	2.25	0.51
1:B:258:GLY:H	1:B:338:GLY:HA3	1.75	0.51
1:A:135:LYS:HD2	1:A:136:SER:N	2.26	0.51
1:A:168:SER:HA	1:A:209:PRO:HD3	1.91	0.51
1:A:95:GLN:HA	1:A:98:LEU:HD12	1.93	0.51
1:A:515:PHE:N	1:A:516:PRO:HD2	2.25	0.51
1:B:100:LEU:HB3	1:B:473:ILE:HG22	1.92	0.50
1:B:104:SER:OG	1:B:473:ILE:O	2.25	0.50
1:A:252:ARG:NE	1:A:335:GLU:HA	2.27	0.50
1:B:202:LEU:O	1:B:241:ARG:NH2	2.41	0.50
1:A:396:PHE:CG	1:A:400:LEU:HD21	2.48	0.49
1:A:89:GLN:O	1:A:93:VAL:HG13	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:323:MET:HG3	1:A:327:LEU:HD12	1.94	0.49
1:B:345:PRO:HA	1:B:368:ALA:HB3	1.94	0.49
1:B:276:ALA:O	1:B:277:HIS:HD2	1.96	0.49
1:A:380:ILE:HG22	1:A:398:GLU:HG2	1.94	0.49
1:A:499:HIS:O	1:A:503:ARG:HG3	2.13	0.48
1:A:254:VAL:HG11	1:A:360:ARG:HG3	1.95	0.48
1:A:252:ARG:HE	1:A:335:GLU:HA	1.79	0.48
1:A:141:PHE:CE1	1:A:153:LEU:HD13	2.50	0.47
1:B:344:MET:HE2	1:B:353:THR:HA	1.96	0.47
1:B:558:TRP:CZ3	1:B:573:ARG:HD2	2.48	0.47
1:A:340:ALA:HA	1:A:363:PRO:HD2	1.96	0.47
1:A:443:ASN:HD21	1:A:516:PRO:HG3	1.78	0.47
1:B:384:ILE:HD12	1:B:403:LEU:HD11	1.96	0.47
1:B:618:LYS:HD2	1:B:618:LYS:N	2.30	0.47
1:A:269:LEU:HB3	1:A:344:MET:HG2	1.96	0.47
1:A:425:ARG:NH2	1:A:507:ASN:O	2.48	0.46
1:A:29:LEU:HD22	1:A:90:ARG:HD3	1.97	0.46
1:B:604:PRO:O	1:B:608:VAL:HG23	2.15	0.46
1:A:181:THR:HG22	1:A:183:THR:H	1.81	0.46
1:A:99:ALA:HA	1:A:146:LEU:HD13	1.96	0.46
1:A:383:VAL:HG23	1:A:398:GLU:HB2	1.97	0.46
1:B:31:ILE:HG23	1:B:45:ILE:HD13	1.98	0.45
1:A:322:ASN:OD1	1:A:322:ASN:N	2.49	0.45
1:A:141:PHE:HE1	1:A:153:LEU:HD13	1.81	0.45
1:A:270:LEU:HD11	1:A:312:PHE:CE2	2.51	0.45
1:B:276:ALA:C	1:B:277:HIS:HD2	2.20	0.45
1:B:477:TYR:O	1:B:480:ARG:HB3	2.17	0.45
1:B:214:HIS:CD2	1:B:214:HIS:H	2.35	0.45
1:B:431:VAL:HB	1:B:461:VAL:HG12	1.99	0.45
1:A:388:ASP:OD2	1:A:552:ARG:NE	2.38	0.45
1:A:29:LEU:HB3	1:A:90:ARG:CZ	2.46	0.45
1:A:540:GLU:HB2	1:A:543:GLU:HG3	1.99	0.45
1:B:515:PHE:CZ	1:B:567:TYR:CD1	3.05	0.45
1:A:204:ASN:ND2	1:A:206:ASN:HB2	2.31	0.44
1:B:441:LYS:HG2	1:B:517:PHE:HB3	1.98	0.44
1:B:527:VAL:HG21	1:B:553:LEU:HD13	1.99	0.44
1:B:396:PHE:CD2	1:B:400:LEU:HD11	2.52	0.44
1:A:42:ILE:HD12	1:A:47:PHE:CE1	2.52	0.44
1:A:3:ASN:HA	1:A:6:LYS:HD3	1.98	0.44
1:B:234:ALA:O	1:B:238:VAL:HG23	2.18	0.44
1:A:446:PHE:HA	1:A:515:PHE:CE1	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:270:LEU:HD11	1:B:312:PHE:CE2	2.53	0.44
1:A:199:LEU:HD12	1:A:202:LEU:HD12	1.99	0.44
1:B:369:LEU:HB3	1:B:372:PRO:HA	2.00	0.44
1:B:396:PHE:CG	1:B:400:LEU:HD21	2.52	0.44
1:A:32:LEU:HB3	1:A:94:TYR:HB3	1.99	0.44
1:A:360:ARG:HD3	1:A:360:ARG:HA	1.81	0.44
1:A:347:ILE:HD12	1:A:353:THR:HB	2.00	0.43
1:B:21:TYR:OH	1:B:78:ASP:OD2	2.30	0.43
1:A:178:PHE:HB2	1:A:439:THR:CB	2.26	0.43
1:B:93:VAL:HG23	1:B:209:PRO:HB3	2.00	0.43
1:B:179:VAL:O	1:B:179:VAL:HG23	2.19	0.43
1:A:217:TYR:O	1:A:232:LYS:HE2	2.18	0.43
1:A:360:ARG:HD2	1:A:362:ALA:O	2.19	0.43
1:B:192:LEU:HD23	1:B:192:LEU:HA	1.89	0.43
1:B:341:ILE:HD13	1:B:364:ILE:HB	2.00	0.43
1:A:340:ALA:HB1	1:A:613:LEU:HD13	2.00	0.43
1:A:510:MET:HB3	1:A:581:ARG:HD2	2.01	0.43
1:B:467:LEU:HD13	1:B:478:VAL:HG11	2.00	0.43
1:A:515:PHE:CB	1:A:536:LYS:HB3	2.49	0.43
1:A:566:GLU:O	1:A:570:ARG:HG2	2.19	0.43
1:B:450:LEU:HB3	1:B:486:LEU:HD11	2.00	0.43
1:B:515:PHE:HB2	1:B:516:PRO:CD	2.49	0.43
1:A:33:SER:HA	1:A:94:TYR:CE1	2.54	0.42
1:B:43:HIS:O	1:B:44:GLU:CB	2.67	0.42
1:B:153:LEU:HD21	1:B:158:MET:CE	2.50	0.42
1:A:139:ILE:HG22	1:A:140:LYS:N	2.34	0.42
1:A:360:ARG:NH1	1:A:365:GLN:HG3	2.34	0.42
1:B:258:GLY:CA	1:B:338:GLY:HA3	2.49	0.42
1:B:588:ILE:HD13	1:B:588:ILE:HA	1.85	0.42
1:A:583:GLU:HG3	1:A:586:ARG:HH21	1.84	0.42
1:B:465:PHE:CE1	1:B:490:ALA:HB1	2.50	0.42
1:B:241:ARG:O	1:B:245:SER:OG	2.37	0.41
1:A:330:ILE:HG22	1:A:361:LEU:HD13	2.02	0.41
1:B:479:GLU:HG3	1:B:483:LYS:HE2	2.02	0.41
1:B:520:THR:H	3:B:701:UDP:PB	2.44	0.41
1:B:540:GLU:HB2	1:B:543:GLU:HG3	2.02	0.41
1:A:12:PHE:CE1	1:A:28:LEU:HB2	2.56	0.41
1:A:249:TRP:NE1	1:A:251:ASP:OD1	2.54	0.41
1:A:277:HIS:O	1:A:281:ARG:HG3	2.20	0.41
1:B:347:ILE:HG21	1:B:367:ILE:HD13	2.03	0.41
1:A:409:PRO:HD2	1:A:602:PRO:HB3	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:515:PHE:HB3	1:A:536:LYS:HB3	2.02	0.41
1:B:344:MET:HB2	1:B:367:ILE:HG22	2.03	0.41
1:B:429:GLU:HG2	1:B:460:LYS:HG3	2.02	0.41
1:A:138:LEU:O	1:A:139:ILE:C	2.58	0.40
1:B:299:LEU:HD22	1:B:316:HIS:HB2	2.03	0.40
1:B:396:PHE:CZ	1:B:541:VAL:HG23	2.56	0.40
1:B:446:PHE:HA	1:B:515:PHE:CD2	2.56	0.40
1:B:603:ARG:N	1:B:604:PRO:CD	2.84	0.40
1:B:608:VAL:O	1:B:612:LYS:HG2	2.20	0.40
1:A:29:LEU:HD22	1:A:90:ARG:CG	2.51	0.40
1:B:447:LEU:HB3	1:B:485:TYR:CD1	2.55	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:414:ALA:O	1:B:118:ASN:ND2[4_455]	2.09	0.11
1:A:429:GLU:OE1	1:B:425:ARG:NH1[1_545]	2.16	0.04

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	604/628 (96%)	590 (98%)	14 (2%)	0	100	100
1	B	590/628 (94%)	575 (98%)	15 (2%)	0	100	100
All	All	1194/1256 (95%)	1165 (98%)	29 (2%)	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	532/552 (96%)	522 (98%)	10 (2%)	57	84
1	B	522/552 (95%)	514 (98%)	8 (2%)	65	87
All	All	1054/1104 (96%)	1036 (98%)	18 (2%)	60	86

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	GLU
1	A	135	LYS
1	A	188	ARG
1	A	303	SER
1	A	357	SER
1	A	438	THR
1	A	499	HIS
1	A	513	ASN
1	A	616	PHE
1	A	618	LYS
1	B	136	SER
1	B	246	GLU
1	B	277	HIS
1	B	285	THR
1	B	501	TYR
1	B	517	PHE
1	B	610	LEU
1	B	618	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	204	ASN
1	A	277	HIS
1	A	427	ASN
1	A	519	ASN

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Mol	Chain	Res	Type
1	A	521	ASN
1	A	579	GLN
1	B	214	HIS
1	B	277	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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
2	UPG	A	701	-	35,38,38	0.96	3 (8%)	53,58,58	1.50	8 (15%)
3	UDP	B	701	-	24,26,26	1.39	4 (16%)	37,40,40	1.93	7 (18%)

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
2	UPG	A	701	-	-	2/23/59/59	0/3/3/3
3	UDP	B	701	-	-	1/16/32/32	0/2/2/2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	701	UDP	C2-N1	3.32	1.43	1.38
3	B	701	UDP	C4-N3	-3.00	1.33	1.38
2	A	701	UPG	C4-N3	-2.53	1.34	1.38
2	A	701	UPG	C2-N3	-2.22	1.34	1.38
3	B	701	UDP	O2-C2	2.17	1.27	1.23
2	A	701	UPG	C5-C4	-2.09	1.39	1.43
3	B	701	UDP	C6-C5	2.06	1.39	1.35

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	701	UDP	N3-C2-N1	5.16	121.73	114.89
2	A	701	UPG	C4-N3-C2	-4.52	120.62	126.58
3	B	701	UDP	O4'-C1'-N1	4.29	118.16	108.36
2	A	701	UPG	N3-C2-N1	4.19	120.45	114.89
3	B	701	UDP	C4-N3-C2	-3.94	121.38	126.58
2	A	701	UPG	PB-O3A-PA	-3.72	120.06	132.83
2	A	701	UPG	C5-C4-N3	3.50	120.08	114.84
3	B	701	UDP	C6-N1-C2	-3.38	116.67	120.99
3	B	701	UDP	C3'-C2'-C1'	3.20	107.50	101.43
2	A	701	UPG	O4-C4-C5	-3.07	119.75	125.16
2	A	701	UPG	C3C-C2C-C1C	2.75	106.65	101.43
3	B	701	UDP	O2-C2-N1	-2.52	119.44	122.79
3	B	701	UDP	C5-C4-N3	2.47	118.54	114.84
2	A	701	UPG	O2-C2-N1	-2.15	119.92	122.79
2	A	701	UPG	O2B-PB-O3B	2.15	115.27	106.78

There are no chirality outliers.

All (3) torsion outliers are listed below:

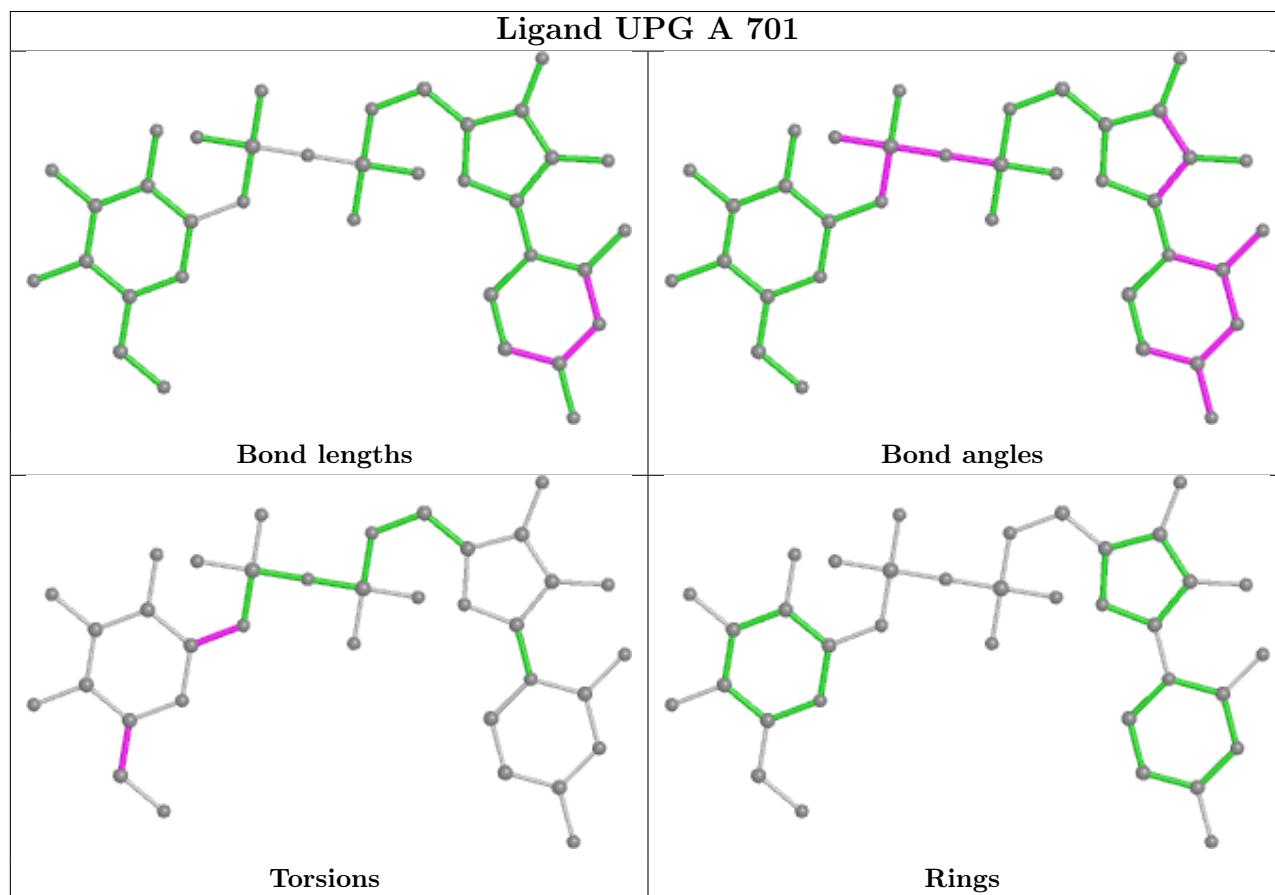
Mol	Chain	Res	Type	Atoms
2	A	701	UPG	O5'-C1'-O3B-PB
2	A	701	UPG	O5'-C5'-C6'-O6'
3	B	701	UDP	C5'-O5'-PA-O1A

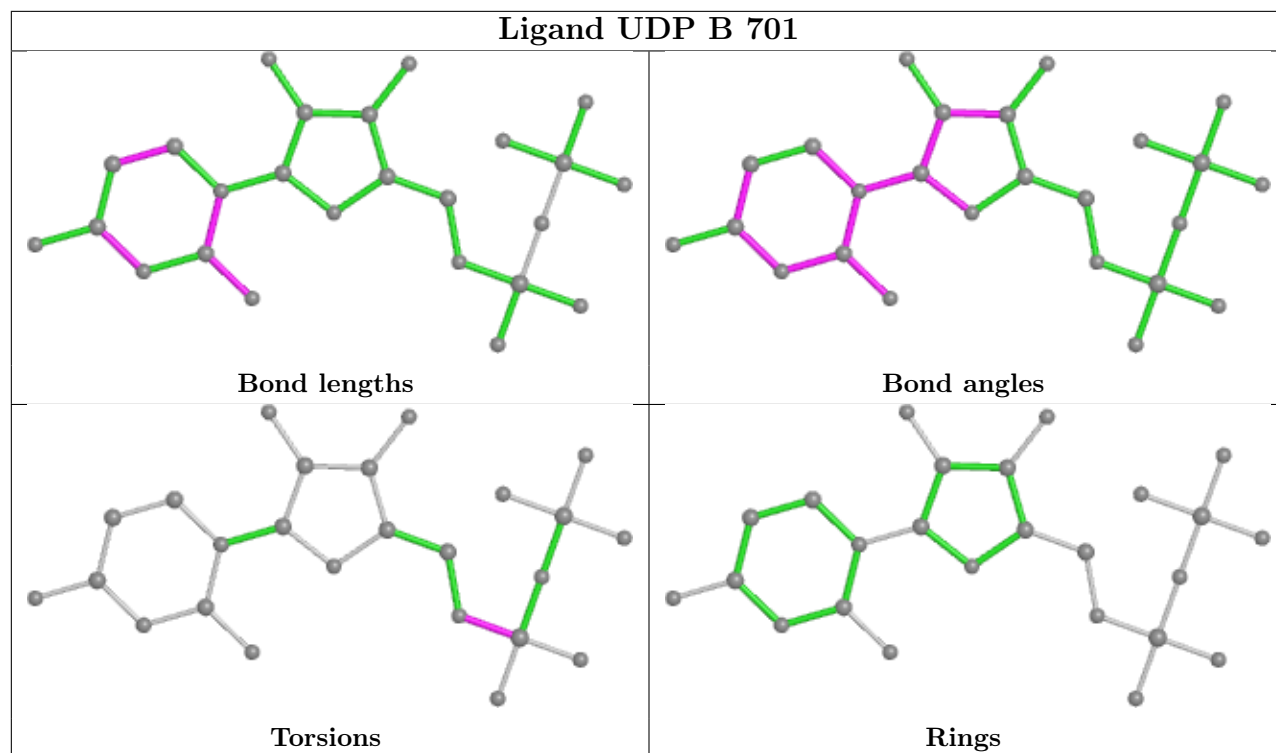
There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	701	UPG	3	0
3	B	701	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	A	608/628 (96%)	0.15	34 (5%) 24 20	28, 45, 70, 104	0
1	B	596/628 (94%)	0.13	22 (3%) 41 37	29, 46, 69, 123	0
All	All	1204/1256 (95%)	0.14	56 (4%) 31 28	28, 45, 69, 123	0

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	487	GLY	7.5
1	A	1	MET	6.2
1	A	133	SER	5.6
1	B	133	SER	4.5
1	B	412	PRO	4.3
1	A	136	SER	4.1
1	A	134	SER	4.1
1	B	132	ASP	4.0
1	B	488	ASP	3.9
1	B	515	PHE	3.8
1	A	120	GLU	3.6
1	A	119	ARG	3.6
1	B	419	LYS	3.4
1	B	420	VAL	3.4
1	B	619	GLU	3.3
1	A	138	LEU	3.2
1	B	429	GLU	3.1
1	A	4	GLU	3.0
1	B	485	TYR	3.0
1	A	135	LYS	3.0
1	A	256	HIS	2.9
1	B	497	PRO	2.9
1	B	616	PHE	2.8
1	A	515	PHE	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	117	TYR	2.7
1	A	591	ASN	2.7
1	A	618	LYS	2.6
1	B	617	LEU	2.6
1	A	334	CYS	2.6
1	A	277	HIS	2.6
1	A	319	ALA	2.5
1	A	161	ILE	2.5
1	A	95	GLN	2.5
1	B	256	HIS	2.5
1	A	139	ILE	2.5
1	A	137	SER	2.4
1	A	619	GLU	2.4
1	A	121	PRO	2.4
1	A	159	TRP	2.3
1	B	430	VAL	2.3
1	A	118	ASN	2.3
1	A	562	ASN	2.3
1	A	116	THR	2.2
1	B	135	LYS	2.2
1	A	318	VAL	2.2
1	B	618	LYS	2.1
1	B	254	VAL	2.1
1	B	6	LYS	2.1
1	A	2	GLU	2.1
1	A	243	ILE	2.1
1	A	114	LEU	2.1
1	A	280	TYR	2.0
1	B	486	LEU	2.0
1	A	90	ARG	2.0
1	A	255	ALA	2.0
1	B	596	THR	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](#)

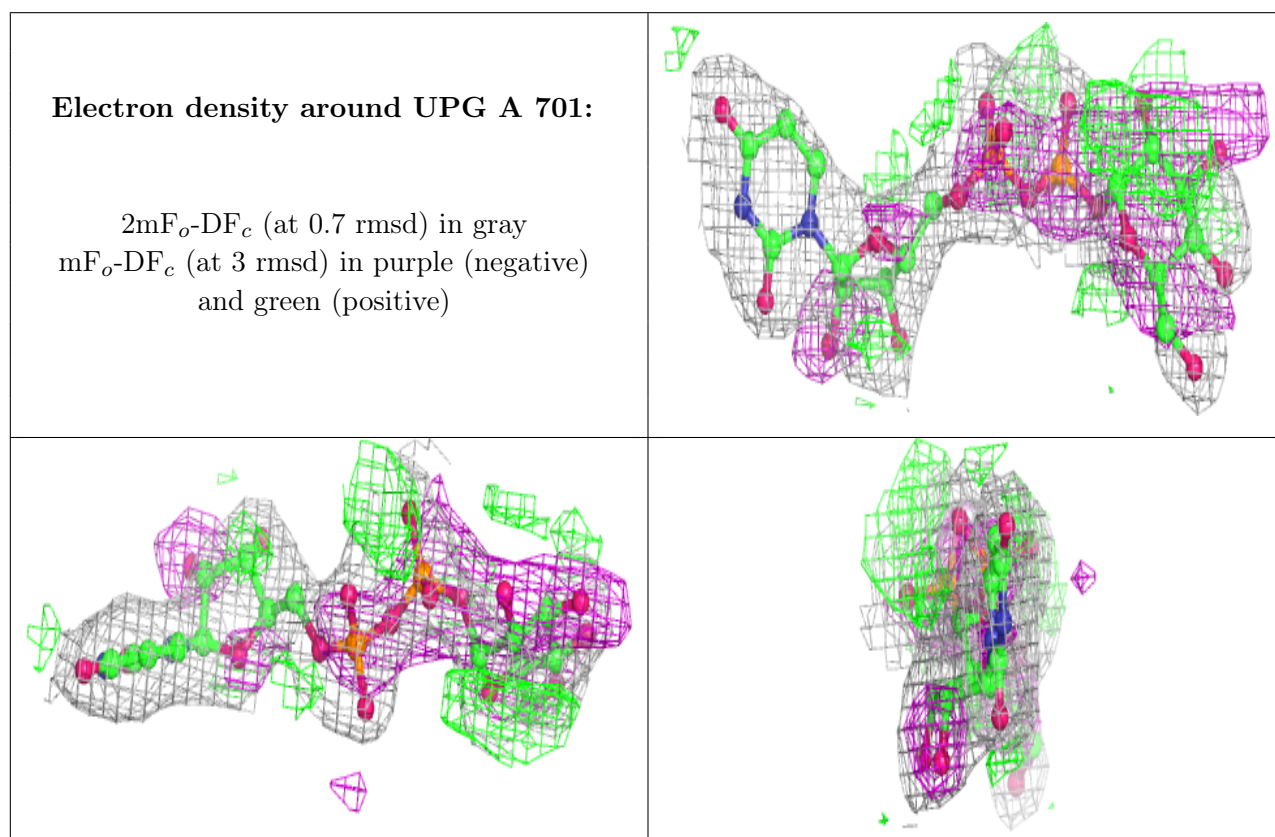
There are no monosaccharides in this entry.

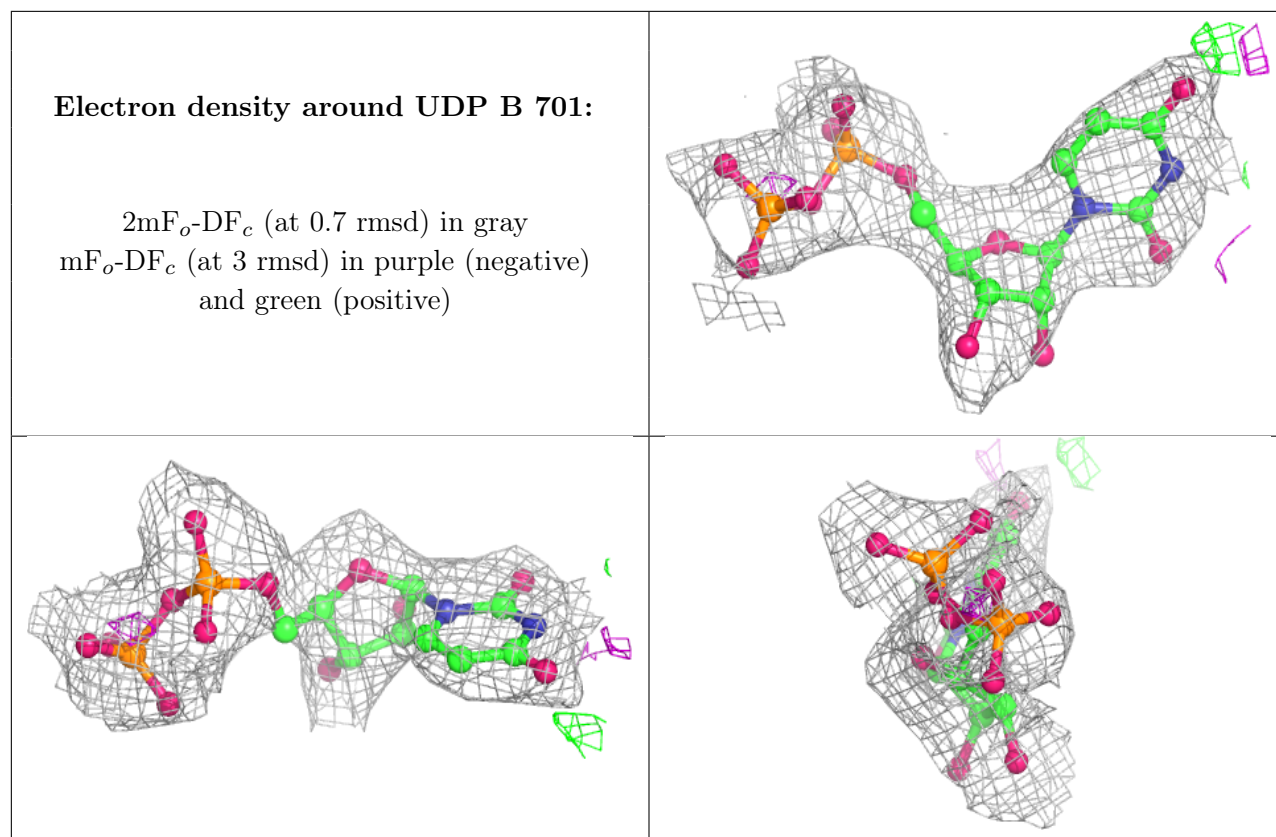
6.4 Ligands [i](#)

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
2	UPG	A	701	36/36	0.81	0.27	20,20,20,20	0
3	UDP	B	701	25/25	0.91	0.17	50,63,71,74	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.