



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 11:32 am BST

PDB ID : 2CVV
Title : Structures of Yeast Ribonucleotide Reductase I
Authors : Xu, H.; Faber, C.; Uchiki, T.; Fairman, J.W.; Racca, J.; Dealwis, C.
Deposited on : 2005-06-14
Resolution : 2.90 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 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 : 2.11
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) : 2.11

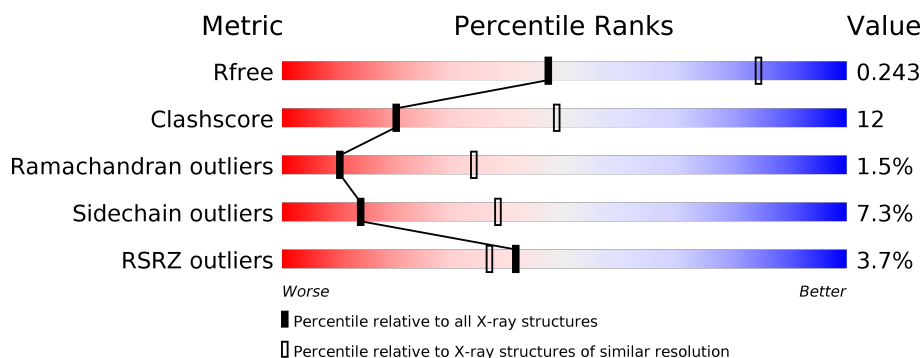
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	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 on 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	888	<div> <div>3%</div> <div> <div></div> <div>50%</div> <div>20%</div> <div>•</div> <div>27%</div> </div> </div>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5403 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.

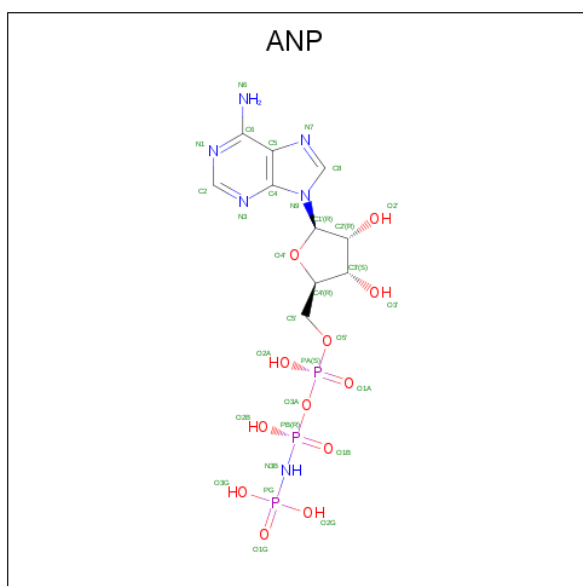
- Molecule 1 is a protein called Ribonucleoside-diphosphate reductase large chain 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	651	Total	C	N	O	S	0	0	0
			5208	3317	883	977	31			

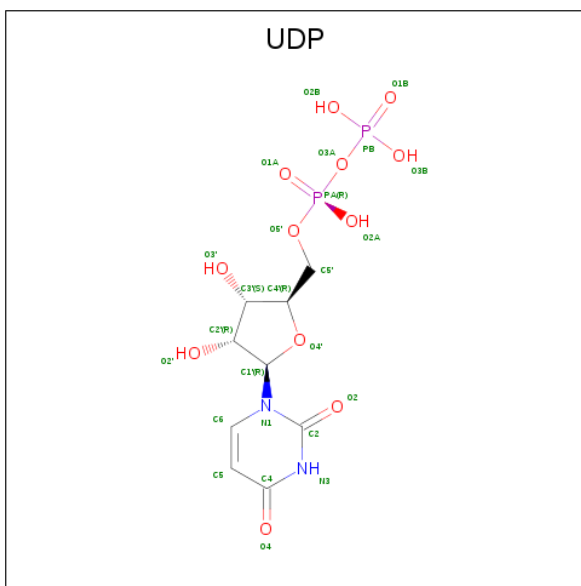
- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (three-letter code: ANP) (formula: $C_{10}H_{17}N_6O_{12}P_3$).



- Molecule 4 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: $\text{C}_9\text{H}_{14}\text{N}_2\text{O}_{12}\text{P}_2$).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	138	Total O 138 138	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	107.87Å 117.59Å 64.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.90 25.99 – 2.60	Depositor EDS
% Data completeness (in resolution range)	97.0 (50.00-2.90) 97.8 (25.99-2.60)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.49 (at 2.60Å)	Xtriage
Refinement program	REFMAC 5.2.0007	Depositor
R, R_{free}	0.177 , 0.245 0.179 , 0.243	Depositor DCC
R_{free} test set	2565 reflections (10.04%)	wwPDB-VP
Wilson B-factor (Å ²)	31.2	Xtriage
Anisotropy	0.688	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 50.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5403	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.32% 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: MG, ANP, UDP

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.65	0/5329	0.85	18/7214 (0.2%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	100	ASP	CB-CG-OD2	7.18	124.76	118.30
1	A	310	ASP	CB-CG-OD2	6.84	124.46	118.30
1	A	233	ASP	CB-CG-OD2	6.66	124.30	118.30
1	A	182	ASP	CB-CG-OD2	6.53	124.18	118.30
1	A	682	ASP	CB-CG-OD2	6.06	123.75	118.30
1	A	138	ASP	CB-CG-OD2	5.99	123.69	118.30
1	A	140	ASP	CB-CG-OD2	5.94	123.65	118.30
1	A	226	ASP	CB-CG-OD2	5.70	123.43	118.30
1	A	586	ASP	CB-CG-OD2	5.61	123.35	118.30
1	A	459	ASP	CB-CG-OD2	5.54	123.29	118.30
1	A	287	ASP	CB-CG-OD2	5.46	123.21	118.30
1	A	438	ASP	CB-CG-OD2	5.22	123.00	118.30
1	A	503	ARG	NE-CZ-NH1	5.15	122.88	120.30
1	A	327	ASP	CB-CG-OD2	5.09	122.88	118.30
1	A	701	ASP	CB-CG-OD2	5.08	122.88	118.30
1	A	129	ASP	CB-CG-OD2	5.06	122.85	118.30
1	A	365	ASP	CB-CG-OD2	5.05	122.84	118.30
1	A	361	ASP	CB-CG-OD2	5.03	122.83	118.30

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	A	5208	0	5135	129	0
2	A	1	0	0	0	0
3	A	31	0	13	0	0
4	A	25	0	11	3	0
5	A	138	0	0	4	0
All	All	5403	0	5159	130	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 12.

All (130) 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:578:LYS:HE2	1:A:578:LYS:HA	1.46	0.95
1:A:505:ILE:HG22	1:A:602:THR:HA	1.51	0.92
1:A:393:LEU:HD22	1:A:724:LEU:HD13	1.48	0.92
1:A:482:ASN:HD22	1:A:599:ASN:HD21	1.18	0.89
1:A:214:GLN:HE22	1:A:216:SER:HB2	1.39	0.88
1:A:272:LEU:O	1:A:276:ILE:HG12	1.77	0.84
1:A:534:GLN:O	1:A:538:THR:HG22	1.82	0.80
1:A:415:ASN:HA	1:A:564:PRO:HG2	1.64	0.77
1:A:428:CYS:SG	4:A:1002:UDP:H3'	2.26	0.76
1:A:482:ASN:ND2	1:A:599:ASN:HD21	1.85	0.74
1:A:662:TYR:CZ	1:A:666:GLN:HG3	2.26	0.71
1:A:140:ASP:OD2	1:A:168:GLN:HG2	1.90	0.70
1:A:692:GLN:O	1:A:696:ILE:HG12	1.91	0.70
1:A:91:THR:HG21	1:A:96:LYS:HG2	1.75	0.69
1:A:288:GLN:HB3	1:A:293:ARG:NH2	2.08	0.68
1:A:106:ASN:OD1	1:A:109:THR:HG22	1.93	0.68
1:A:602:THR:N	1:A:707:ASP:OD2	2.29	0.65
1:A:693:LYS:H	1:A:693:LYS:HD2	1.62	0.64
1:A:109:THR:HG23	1:A:111:LYS:H	1.63	0.64
1:A:717:ARG:O	1:A:719:PRO:HD3	1.97	0.63
1:A:534:GLN:O	1:A:538:THR:CG2	2.47	0.63
1:A:456:THR:HA	1:A:463:SER:HA	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:475:LYS:HD2	1:A:546:ALA:HB2	1.80	0.63
1:A:481:LEU:CB	1:A:505:ILE:HG12	2.30	0.61
1:A:571:GLN:HA	1:A:571:GLN:HE21	1.65	0.61
1:A:142:GLN:HB3	1:A:196:LYS:HE2	1.82	0.61
1:A:393:LEU:CD2	1:A:724:LEU:HD13	2.29	0.60
1:A:661:GLN:HA	1:A:664:ILE:HD12	1.83	0.60
1:A:665:THR:HG22	1:A:666:GLN:HE21	1.67	0.59
1:A:225:GLU:HG3	5:A:2133:HOH:O	2.03	0.59
1:A:317:ASN:OD1	1:A:326:ARG:NH1	2.36	0.59
1:A:571:GLN:HA	1:A:571:GLN:NE2	2.19	0.58
1:A:373:ARG:HG2	1:A:377:GLU:OE2	2.04	0.58
1:A:481:LEU:HB3	1:A:505:ILE:HG12	1.85	0.58
1:A:165:GLU:HG2	1:A:169:HIS:HB2	1.86	0.57
1:A:214:GLN:HE21	1:A:488:ASN:HD21	1.53	0.56
1:A:220:LEU:HD22	1:A:426:ASN:HB3	1.88	0.55
1:A:538:THR:HB	1:A:583:TRP:NE1	2.22	0.55
1:A:713:ASN:ND2	1:A:742:TYR:H	2.05	0.54
1:A:446:ALA:HB3	1:A:477:VAL:CG1	2.37	0.54
1:A:665:THR:HG22	1:A:666:GLN:NE2	2.22	0.54
1:A:306:ALA:HA	1:A:350:LEU:HB3	1.89	0.54
1:A:146:PHE:CZ	1:A:640:VAL:HG21	2.43	0.54
1:A:578:LYS:HA	1:A:578:LYS:CE	2.31	0.54
1:A:270:ASN:HB3	1:A:274:PRO:HG2	1.90	0.54
1:A:606:MET:HB2	1:A:607:PRO:CD	2.38	0.54
1:A:691:SER:HB2	1:A:693:LYS:HD3	1.89	0.54
1:A:475:LYS:O	1:A:479:ARG:HG3	2.09	0.53
1:A:214:GLN:HE22	1:A:216:SER:CB	2.15	0.53
1:A:723:LYS:HG2	5:A:2115:HOH:O	2.08	0.53
1:A:447:SER:HB3	1:A:606:MET:HE3	1.90	0.52
1:A:560:PHE:CZ	1:A:596:GLY:HA2	2.45	0.52
1:A:655:TRP:CZ2	1:A:660:LYS:HG2	2.45	0.52
1:A:538:THR:HB	1:A:583:TRP:HE1	1.75	0.51
1:A:436:ALA:HB1	1:A:437:PRO:CD	2.41	0.50
1:A:288:GLN:O	1:A:291:ASN:ND2	2.45	0.49
1:A:654:ILE:O	1:A:659:MET:HG3	2.13	0.49
1:A:304:TRP:CZ2	1:A:359:LEU:HB3	2.48	0.49
1:A:481:LEU:HB2	1:A:505:ILE:HG12	1.94	0.48
1:A:450:LEU:HD22	1:A:535:ILE:HG21	1.95	0.48
1:A:693:LYS:H	1:A:693:LYS:CD	2.26	0.48
1:A:94:PHE:HB3	1:A:132:ASN:OD1	2.14	0.48
1:A:298:ALA:HB2	1:A:427:LEU:HD12	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:ARG:HD3	1:A:194:SER:HB2	1.96	0.48
1:A:168:GLN:NE2	1:A:190:TYR:OH	2.43	0.48
1:A:482:ASN:HD21	1:A:503:ARG:HH11	1.61	0.48
1:A:334:ILE:HD12	1:A:404:VAL:HG13	1.97	0.47
1:A:251:HIS:HB3	1:A:424:SER:HB3	1.96	0.47
1:A:168:GLN:HE22	1:A:194:SER:HB3	1.80	0.47
1:A:220:LEU:N	1:A:220:LEU:HD23	2.30	0.46
1:A:673:LEU:HD22	1:A:676:VAL:HG23	1.98	0.46
1:A:625:THR:HA	1:A:687:VAL:HG12	1.97	0.46
1:A:288:GLN:HB3	1:A:293:ARG:HH21	1.77	0.46
1:A:482:ASN:ND2	1:A:503:ARG:HH11	2.13	0.46
1:A:520:ARG:HH22	1:A:648:ASP:CG	2.18	0.46
1:A:606:MET:HE3	1:A:608:THR:HG22	1.98	0.46
1:A:740:MET:SD	1:A:743:LEU:HB2	2.56	0.46
1:A:311:PHE:HA	1:A:314:ILE:HD13	1.98	0.45
1:A:557:TYR:HB3	1:A:598:ARG:O	2.17	0.45
1:A:104:TYR:CD1	1:A:159:ILE:HG23	2.52	0.45
1:A:444:ASN:C	1:A:445:LEU:HD12	2.37	0.45
1:A:458:GLU:OE1	1:A:462:THR:HB	2.16	0.45
1:A:483:ARG:HH22	1:A:487:ARG:HD2	1.82	0.45
1:A:512:LEU:HB2	1:A:621:PHE:HA	2.00	0.44
1:A:288:GLN:C	1:A:290:GLY:H	2.20	0.44
1:A:662:TYR:CE1	1:A:666:GLN:HG3	2.51	0.44
1:A:181:ARG:O	1:A:183:ILE:HD12	2.18	0.44
1:A:412:ARG:O	1:A:417:LYS:NZ	2.37	0.44
1:A:428:CYS:HB2	1:A:430:GLU:OE2	2.17	0.44
1:A:692:GLN:NE2	1:A:715:PHE:H	2.15	0.44
1:A:604:ALA:HB2	1:A:708:GLN:HB2	2.00	0.43
1:A:136:VAL:HB	1:A:139:ARG:HD2	2.00	0.43
1:A:671:GLN:O	1:A:672:GLY:C	2.56	0.43
1:A:507:LEU:HD12	1:A:602:THR:O	2.19	0.43
1:A:120:VAL:HG21	1:A:209:GLY:HA2	2.01	0.43
1:A:436:ALA:HB1	1:A:437:PRO:HD2	2.01	0.43
1:A:601:LEU:HA	1:A:707:ASP:OD2	2.19	0.42
1:A:164:ALA:HB3	5:A:2071:HOH:O	2.18	0.42
1:A:109:THR:OG1	1:A:111:LYS:HG3	2.19	0.42
1:A:442:VAL:HG12	1:A:443:CYS:N	2.34	0.42
1:A:401:THR:HB	1:A:402:PRO:HA	2.02	0.42
4:A:1002:UDP:H6	4:A:1002:UDP:O5'	2.02	0.42
1:A:446:ALA:HB3	1:A:477:VAL:HG11	2.00	0.42
1:A:577:GLN:HB3	5:A:2106:HOH:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:509:VAL:O	1:A:620:CYS:HA	2.20	0.42
1:A:413:LYS:HE3	1:A:575:TRP:CE2	2.55	0.42
1:A:518:LEU:HA	1:A:518:LEU:HD12	1.94	0.41
1:A:628:MET:HE2	1:A:664:ILE:HG12	2.02	0.41
1:A:131:LEU:HD12	1:A:131:LEU:HA	1.82	0.41
1:A:450:LEU:HD22	1:A:535:ILE:HD13	2.01	0.41
1:A:455:GLU:C	1:A:457:SER:H	2.24	0.41
1:A:127:ASN:HB2	1:A:131:LEU:HD22	2.02	0.41
1:A:91:THR:HG21	1:A:96:LYS:CG	2.47	0.41
1:A:293:ARG:HA	1:A:294:PRO:HD3	1.82	0.41
1:A:482:ASN:ND2	1:A:503:ARG:NH1	2.69	0.41
1:A:426:ASN:ND2	4:A:1002:UDP:O3'	2.53	0.41
1:A:256:ARG:HG2	1:A:260:SER:HB2	2.03	0.41
1:A:659:MET:CE	1:A:673:LEU:HD11	2.51	0.41
1:A:673:LEU:HD22	1:A:676:VAL:CG2	2.50	0.41
1:A:170:LEU:C	1:A:170:LEU:HD23	2.41	0.41
1:A:259:GLY:HA2	1:A:268:THR:HG23	2.02	0.41
1:A:484:VAL:O	1:A:488:ASN:HB2	2.21	0.41
1:A:220:LEU:HD21	1:A:431:ILE:HG21	2.01	0.41
1:A:713:ASN:HD22	1:A:742:TYR:H	1.69	0.41
1:A:702:ARG:O	1:A:706:ILE:HG13	2.20	0.41
1:A:588:LEU:O	1:A:592:ILE:HG13	2.21	0.41
1:A:537:GLU:OE2	1:A:582:MET:HB3	2.21	0.40
1:A:713:ASN:ND2	1:A:742:TYR:HB2	2.36	0.40
1:A:578:LYS:HE2	1:A:578:LYS:CA	2.32	0.40
1:A:656:ASP:OD1	1:A:658:GLY:N	2.51	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	647/888 (73%)	591 (91%)	46 (7%)	10 (2%)	10	34

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	296	ALA
1	A	459	ASP
1	A	460	GLY
1	A	629	TYR
1	A	639	GLN
1	A	707	ASP
1	A	256	ARG
1	A	314	ILE
1	A	741	TYR
1	A	161	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	565/761 (74%)	524 (93%)	41 (7%)	14	38

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	119	ASP
1	A	126	GLU
1	A	131	LEU
1	A	146	PHE
1	A	162	GLN
1	A	176	LEU
1	A	183	ILE
1	A	195	LEU
1	A	212	LYS
1	A	214	GLN
1	A	220	LEU

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Mol	Chain	Res	Type
1	A	243	LYS
1	A	268	THR
1	A	292	LYS
1	A	293	ARG
1	A	301	LEU
1	A	314	ILE
1	A	321	GLU
1	A	326	ARG
1	A	337	LEU
1	A	359	LEU
1	A	388	LEU
1	A	443	CYS
1	A	456	THR
1	A	459	ASP
1	A	462	THR
1	A	503	ARG
1	A	505	ILE
1	A	512	LEU
1	A	518	LEU
1	A	520	ARG
1	A	530	LEU
1	A	538	THR
1	A	606	MET
1	A	610	SER
1	A	638	PHE
1	A	639	GLN
1	A	659	MET
1	A	712	LEU
1	A	720	THR
1	A	746	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	168	GLN
1	A	214	GLN
1	A	444	ASN
1	A	482	ASN
1	A	613	GLN
1	A	666	GLN
1	A	692	GLN
1	A	710	HIS

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Mol	Chain	Res	Type
1	A	713	ASN
1	A	746	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	ANP	A	1001	2	29,33,33	1.37	4 (13%)	31,52,52	2.42	13 (41%)
4	UDP	A	1002	-	20,26,26	1.34	1 (5%)	25,40,40	1.01	1 (4%)

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	ANP	A	1001	2	-	1/14/38/38	0/3/3/3
4	UDP	A	1002	-	-	6/14/32/32	0/2/2/2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1002	UDP	C4-N3	4.14	1.40	1.33
3	A	1001	ANP	C5-C4	3.35	1.49	1.40
3	A	1001	ANP	PG-O1G	3.16	1.51	1.46
3	A	1001	ANP	PB-O1B	2.68	1.50	1.46
3	A	1001	ANP	PB-O2B	-2.18	1.50	1.56

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1001	ANP	O1G-PG-N3B	-6.86	101.67	111.77
3	A	1001	ANP	C3'-C2'-C1'	4.88	108.32	100.98
3	A	1001	ANP	O1B-PB-N3B	-3.99	105.90	111.77
3	A	1001	ANP	O2B-PB-O1B	3.65	117.57	109.92
3	A	1001	ANP	O3A-PB-N3B	-3.30	97.43	106.59
3	A	1001	ANP	C4-C5-N7	-3.28	105.98	109.40
3	A	1001	ANP	O4'-C1'-C2'	-2.70	102.98	106.93
3	A	1001	ANP	N3-C2-N1	-2.63	124.56	128.68
3	A	1001	ANP	PA-O3A-PB	-2.41	124.13	132.62
3	A	1001	ANP	C2-N1-C6	2.38	122.82	118.75
3	A	1001	ANP	O2A-PA-O1A	2.35	123.86	112.24
4	A	1002	UDP	C6-N1-C2	-2.18	117.74	121.20
3	A	1001	ANP	C5-C6-N6	2.03	123.44	120.35
3	A	1001	ANP	O2B-PB-O3A	2.00	111.32	104.64

There are no chirality outliers.

All (7) torsion outliers are listed below:

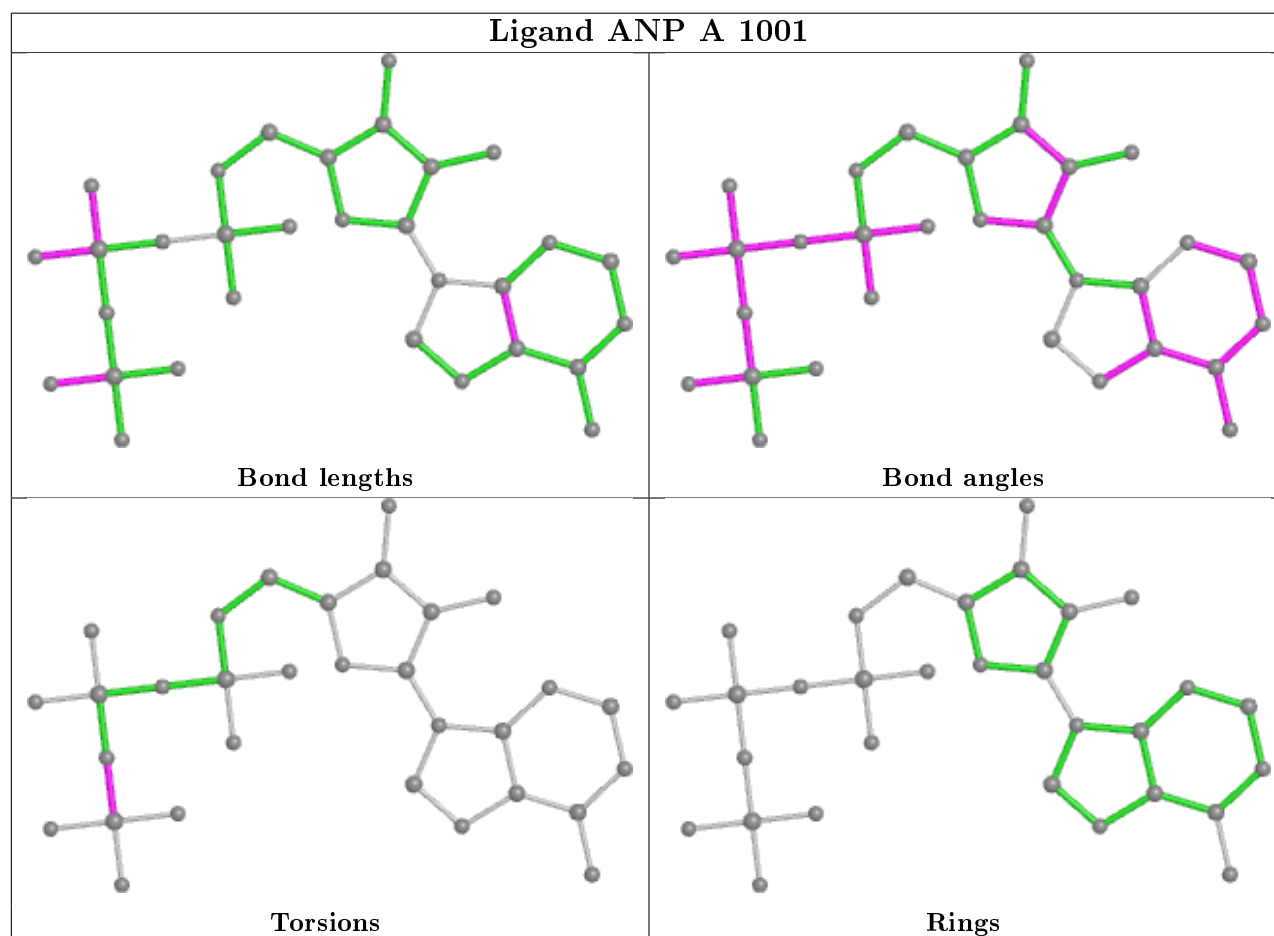
Mol	Chain	Res	Type	Atoms
3	A	1001	ANP	PB-N3B-PG-O1G
4	A	1002	UDP	C5'-O5'-PA-O1A
4	A	1002	UDP	C5'-O5'-PA-O2A
4	A	1002	UDP	O4'-C4'-C5'-O5'
4	A	1002	UDP	C3'-C4'-C5'-O5'
4	A	1002	UDP	PB-O3A-PA-O1A
4	A	1002	UDP	C5'-O5'-PA-O3A

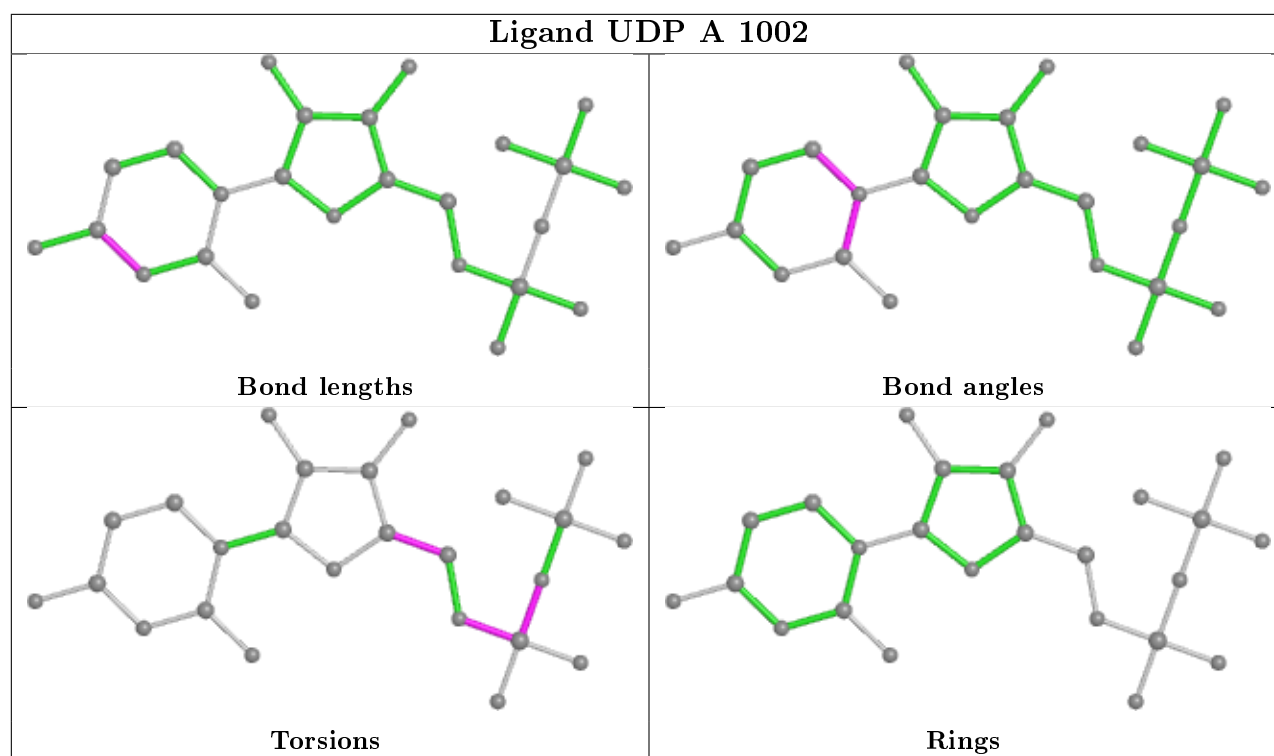
There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1002	UDP	3	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	651/888 (73%)	-0.34	24 (3%)	41 37	12, 24, 57, 69	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	458	GLU	5.2
1	A	629	TYR	5.0
1	A	457	SER	5.0
1	A	294	PRO	4.8
1	A	89	GLN	3.9
1	A	459	ASP	3.9
1	A	292	LYS	3.8
1	A	718	ALA	3.6
1	A	638	PHE	3.5
1	A	319	GLY	3.4
1	A	460	GLY	3.1
1	A	746	GLN	2.8
1	A	293	ARG	2.8
1	A	456	THR	2.7
1	A	461	LYS	2.7
1	A	320	LYS	2.6
1	A	317	ASN	2.5
1	A	91	THR	2.4
1	A	323	ILE	2.3
1	A	90	THR	2.3
1	A	717	ARG	2.3
1	A	462	THR	2.2
1	A	667	ASN	2.2
1	A	661	GLN	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](#)

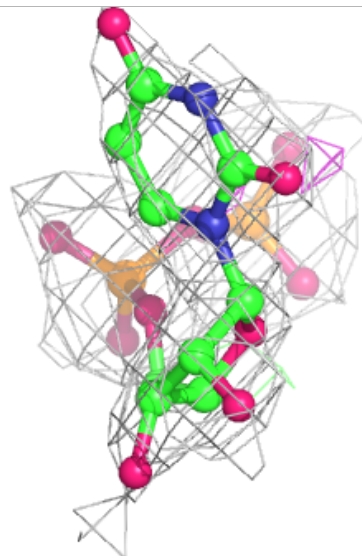
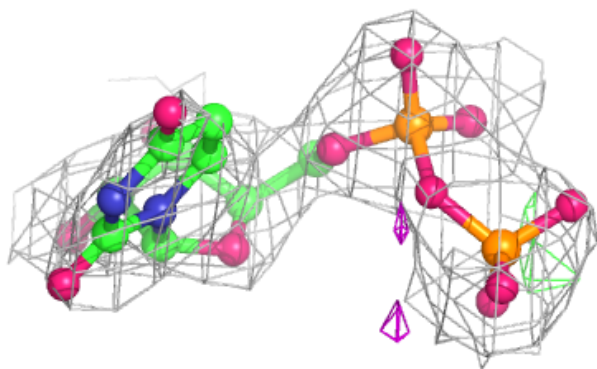
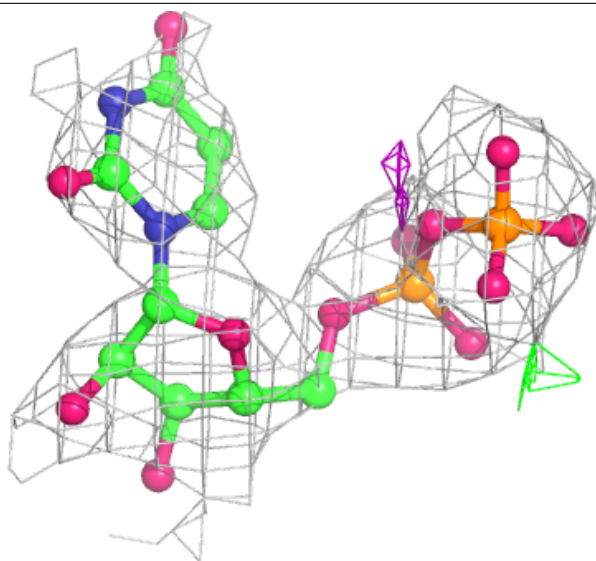
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	UDP	A	1002	25/25	0.93	0.26	40,54,58,58	0
3	ANP	A	1001	31/31	0.97	0.10	14,18,20,26	0
2	MG	A	2001	1/1	0.97	0.05	16,16,16,16	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.

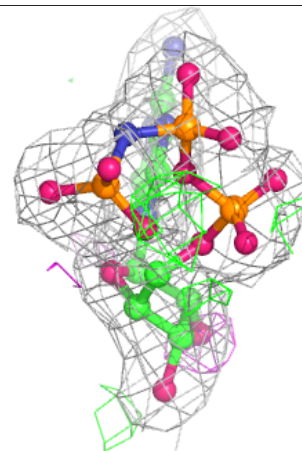
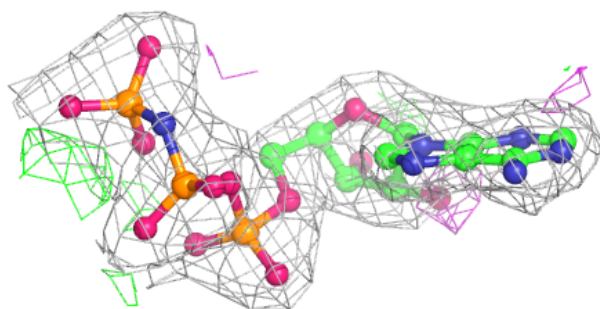
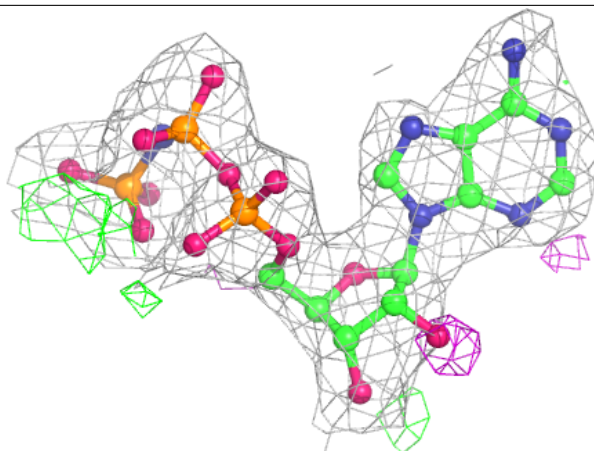
Electron density around UDP A 1002:

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



Electron density around ANP A 1001:

$2mF_o-DF_c$ (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.