



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 07:10 am GMT

PDB ID : 2CVX  
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 : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : recalc28949

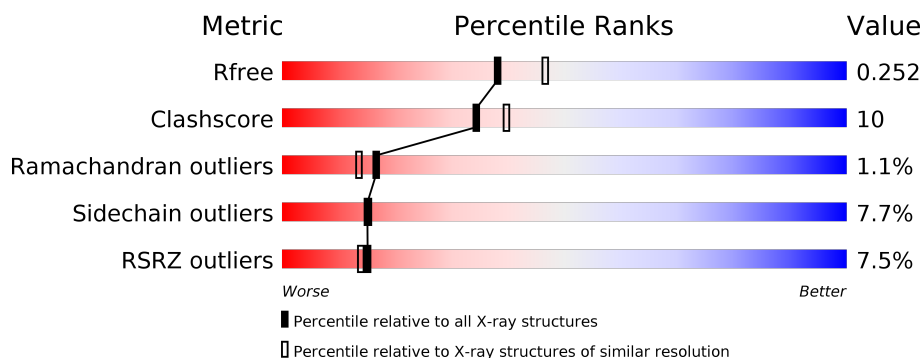
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	100719	4002 (2.20-2.20)
Clashscore	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)
RSRZ outliers	101464	4033 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	888	

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5580 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.

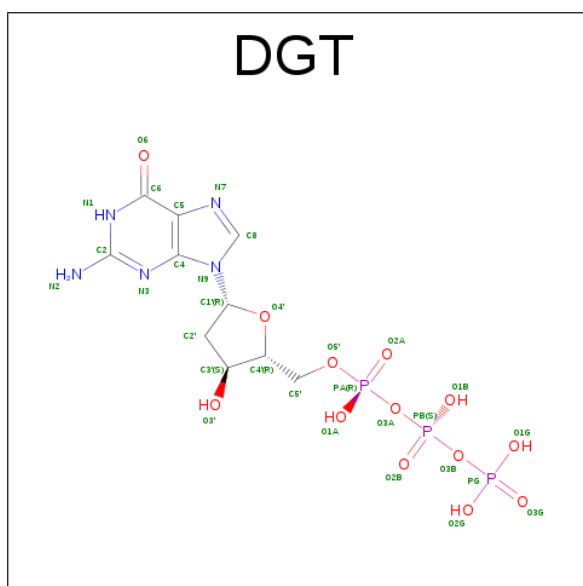
- Molecule 1 is a protein called Ribonucleoside-diphosphate reductase large chain 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	664	Total	C	N	O	S	0	0	0
			5306	3379	903	993	31			

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

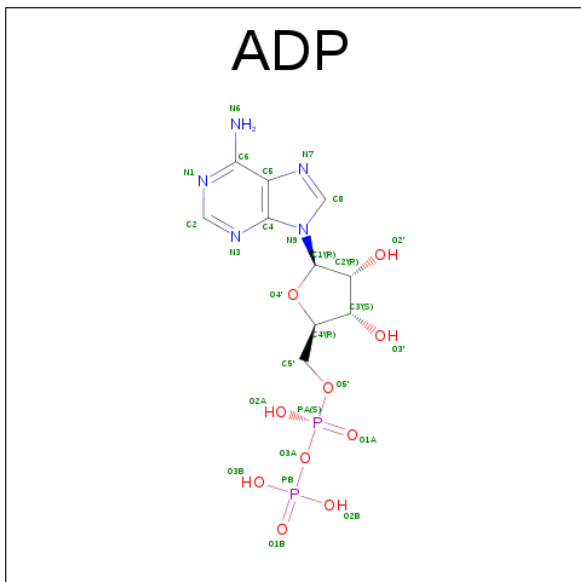
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is 2'-DEOXYGUANOSINE-5'-TRIPHOSPHATE (three-letter code: DGT) (formula: C<sub>10</sub>H<sub>16</sub>N<sub>5</sub>O<sub>13</sub>P<sub>3</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula:  $C_{10}H_{15}N_5O_{10}P_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	215	Total	O	0	0
			215	215		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	107.69Å 117.28Å 65.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.20 33.86 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.2 (50.00-2.20) 99.2 (33.86-2.20)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.61 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.2.0007	Depositor
R, $R_{free}$	0.210 , 0.254 0.209 , 0.252	Depositor DCC
$R_{free}$ test set	4224 reflections (11.09%)	DCC
Wilson B-factor (Å <sup>2</sup> )	31.2	Xtriage
Anisotropy	1.067	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 44.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	5580	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.97% of the height of the origin peak. No significant pseudotranslation is detected.*

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

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

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, ADP, DGT

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	0.59	0/5428	0.84	17/7348 (0.2%)

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	A	0	1

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	126	GLU	OE1-CD-OE2	-10.99	110.11	123.30
1	A	503	ARG	NE-CZ-NH2	-9.05	115.78	120.30
1	A	182	ASP	CB-CG-OD2	7.00	124.60	118.30
1	A	707	ASP	CB-CG-OD2	6.86	124.47	118.30
1	A	138	ASP	CB-CG-OD2	6.58	124.23	118.30
1	A	327	ASP	CB-CG-OD2	6.38	124.04	118.30
1	A	287	ASP	CB-CG-OD2	6.27	123.94	118.30
1	A	503	ARG	NE-CZ-NH1	6.16	123.38	120.30
1	A	438	ASP	CB-CG-OD2	6.07	123.76	118.30
1	A	365	ASP	CB-CG-OD2	5.80	123.52	118.30
1	A	586	ASP	CB-CG-OD2	5.78	123.50	118.30
1	A	100	ASP	CB-CG-OD2	5.67	123.40	118.30
1	A	459	ASP	CB-CG-OD2	5.56	123.30	118.30
1	A	648	ASP	CB-CG-OD2	5.51	123.26	118.30
1	A	170	LEU	CA-CB-CG	5.35	127.60	115.30
1	A	140	ASP	CB-CG-OD2	5.04	122.84	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	591	ASP	CB-CG-OD2	5.01	122.81	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	126	GLU	Sidechain

## 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	5306	0	5245	101	0
2	A	1	0	0	0	0
3	A	31	0	12	1	0
4	A	27	0	12	1	0
5	A	215	0	0	3	0
All	All	5580	0	5269	101	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 10.

All (101) 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:106:ASN:OD1	1:A:109:THR:HG22	1.55	1.06
1:A:534:GLN:O	1:A:538:THR:HG22	1.62	0.97
1:A:522:PRO:HG2	1:A:525:SER:HB3	1.44	0.96
1:A:538:THR:HB	1:A:583:TRP:HE1	1.41	0.83
1:A:538:THR:HB	1:A:583:TRP:NE1	1.96	0.79
1:A:686:THR:HG22	1:A:688:TRP:H	1.52	0.75
1:A:505:ILE:HG22	1:A:602:THR:HA	1.70	0.73
1:A:445:LEU:HD22	1:A:506:ALA:HB3	1.70	0.73
1:A:170:LEU:HD22	1:A:173:ARG:NH2	2.04	0.72
1:A:481:LEU:HB3	1:A:505:ILE:HG13	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:457:SER:HB3	1:A:462:THR:HG23	1.74	0.70
1:A:717:ARG:O	1:A:719:PRO:HD3	1.91	0.70
1:A:319:GLY:O	1:A:320:LYS:HB2	1.92	0.70
1:A:502:HIS:ND1	1:A:559:THR:HG21	2.09	0.67
1:A:485:ILE:HD11	1:A:505:ILE:HD13	1.77	0.67
1:A:508:GLY:HA3	1:A:606:MET:HE1	1.78	0.65
1:A:214:GLN:HE22	1:A:216:SER:HB2	1.61	0.65
1:A:334:ILE:HD12	1:A:404:VAL:HG13	1.80	0.64
1:A:520:ARG:HH22	1:A:648:ASP:CG	2.01	0.64
1:A:447:SER:HB3	1:A:606:MET:HE3	1.79	0.64
1:A:745:THR:O	1:A:746:GLN:HB2	1.99	0.63
1:A:481:LEU:HB3	1:A:505:ILE:CG1	2.30	0.61
1:A:251:HIS:HD2	5:A:2033:HOH:O	1.84	0.61
1:A:662:TYR:CD1	1:A:662:TYR:O	2.54	0.61
1:A:520:ARG:NH2	1:A:648:ASP:OD2	2.33	0.59
1:A:214:GLN:NE2	5:A:2173:HOH:O	2.31	0.59
1:A:428:CYS:SG	4:A:1002:ADP:H3'	2.44	0.57
1:A:686:THR:HG21	1:A:688:TRP:HD1	1.69	0.56
1:A:506:ALA:HB1	1:A:604:ALA:HB3	1.87	0.56
1:A:692:GLN:HB2	1:A:727:MET:SD	2.47	0.55
1:A:401:THR:HB	1:A:402:PRO:HA	1.89	0.55
1:A:589:ARG:O	1:A:593:MET:HG3	2.06	0.55
1:A:145:TYR:N	1:A:145:TYR:CD2	2.72	0.55
1:A:447:SER:HB3	1:A:606:MET:CE	2.37	0.54
1:A:683:LEU:HD23	1:A:684:TYR:CZ	2.43	0.53
1:A:383:ILE:HG13	1:A:384:LYS:H	1.75	0.52
1:A:109:THR:HG23	1:A:111:LYS:HB2	1.89	0.52
1:A:314:ILE:HG12	1:A:325:ALA:HB3	1.91	0.52
1:A:320:LYS:O	1:A:323:ILE:HG22	2.10	0.52
1:A:485:ILE:HD11	1:A:505:ILE:CD1	2.38	0.51
1:A:82:ALA:HA	1:A:85:ASN:HD22	1.75	0.51
1:A:606:MET:HE3	1:A:608:THR:HG22	1.92	0.51
1:A:332:LEU:HD11	1:A:392:ILE:HD12	1.91	0.51
1:A:610:SER:O	1:A:613:GLN:HB3	2.11	0.50
1:A:488:ASN:ND2	5:A:2173:HOH:O	2.31	0.50
1:A:504:PRO:O	1:A:505:ILE:HD12	2.12	0.50
1:A:388:LEU:HD22	1:A:392:ILE:HD11	1.93	0.50
1:A:560:PHE:CZ	1:A:596:GLY:HA2	2.47	0.49
1:A:356:ALA:HB1	1:A:374:TYR:CD1	2.47	0.49
1:A:86:LEU:HD22	1:A:148:PHE:HE1	1.78	0.48
1:A:388:LEU:O	1:A:392:ILE:HG12	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:393:LEU:HD22	1:A:724:LEU:HB3	1.95	0.48
1:A:251:HIS:HB3	1:A:424:SER:HB3	1.96	0.48
1:A:383:ILE:CG1	1:A:384:LYS:H	2.26	0.47
1:A:714:LEU:HD22	1:A:740:MET:HE3	1.96	0.47
1:A:126:GLU:CD	1:A:181:ARG:HH12	2.18	0.47
1:A:713:ASN:ND2	1:A:742:TYR:HB2	2.29	0.47
1:A:519:LEU:O	1:A:520:ARG:HB2	2.15	0.47
1:A:242:SER:O	1:A:288:GLN:HA	2.15	0.47
1:A:220:LEU:HD21	1:A:426:ASN:HB3	1.97	0.47
1:A:740:MET:SD	1:A:743:LEU:HB2	2.54	0.47
1:A:297:PHE:HB2	1:A:328:LEU:HD22	1.96	0.47
1:A:106:ASN:OD1	1:A:109:THR:CG2	2.44	0.46
1:A:606:MET:HE3	1:A:608:THR:CG2	2.46	0.45
1:A:220:LEU:CD2	1:A:426:ASN:HB3	2.47	0.45
1:A:86:LEU:HD22	1:A:148:PHE:CE1	2.51	0.45
1:A:594:LYS:O	1:A:594:LYS:HE3	2.17	0.45
1:A:383:ILE:HG13	1:A:384:LYS:N	2.32	0.45
1:A:475:LYS:O	1:A:479:ARG:HG3	2.16	0.45
1:A:179:HIS:ND1	1:A:483:ARG:NH1	2.65	0.45
1:A:692:GLN:NE2	1:A:715:PHE:H	2.15	0.44
1:A:273:ILE:HG21	1:A:323:ILE:HA	1.99	0.44
1:A:210:THR:HB	1:A:211:PRO:CD	2.48	0.44
1:A:686:THR:CG2	1:A:688:TRP:HD1	2.31	0.44
1:A:273:ILE:HB	1:A:274:PRO:HD3	2.00	0.44
1:A:338:PHE:O	1:A:342:VAL:HG23	2.17	0.43
1:A:79:ALA:C	1:A:81:ILE:H	2.21	0.43
1:A:109:THR:HG23	1:A:111:LYS:H	1.83	0.43
1:A:264:GLY:HA3	3:A:1001:DGT:O2B	2.18	0.43
1:A:719:PRO:HG3	1:A:745:THR:OG1	2.19	0.43
1:A:662:TYR:O	1:A:662:TYR:HD1	2.00	0.43
1:A:288:GLN:HG3	1:A:293:ARG:O	2.19	0.43
1:A:413:LYS:NZ	1:A:735:GLY:O	2.52	0.42
1:A:432:VAL:CG1	1:A:708:GLN:HA	2.48	0.42
1:A:76:THR:HG23	1:A:77:LEU:H	1.83	0.42
1:A:248:ILE:HD12	1:A:297:PHE:CE2	2.54	0.42
1:A:436:ALA:HB1	1:A:437:PRO:HD2	2.01	0.42
1:A:140:ASP:OD2	1:A:168:GLN:HG2	2.20	0.42
1:A:413:LYS:HE2	1:A:575:TRP:CE2	2.55	0.42
1:A:715:PHE:CE2	1:A:744:ARG:HG3	2.55	0.42
1:A:126:GLU:OE1	1:A:181:ARG:NH1	2.53	0.42
1:A:383:ILE:CG1	1:A:384:LYS:N	2.83	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:104:TYR:CD1	1:A:159:ILE:HG23	2.55	0.42
1:A:119:ASP:O	1:A:123:ILE:HG13	2.19	0.42
1:A:178:ILE:O	1:A:487:ARG:HD3	2.20	0.41
1:A:524:ASP:HB3	1:A:694:THR:HG21	2.01	0.41
1:A:557:TYR:CD1	1:A:559:THR:HG22	2.56	0.41
1:A:714:LEU:HD22	1:A:740:MET:CE	2.51	0.41
1:A:450:LEU:N	1:A:451:PRO:CD	2.84	0.41
1:A:120:VAL:O	1:A:124:VAL:HG23	2.21	0.40
1:A:128:LYS:O	1:A:132:ASN:HB2	2.21	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

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

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	660/888 (74%)	624 (94%)	29 (4%)	7 (1%)	17	13

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	320	LYS
1	A	458	GLU
1	A	459	ASP
1	A	80	ARG
1	A	674	PRO
1	A	741	TYR
1	A	620	CYS

### 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	575/761 (76%)	531 (92%)	44 (8%)	15	15

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	145	TYR
1	A	154	SER
1	A	170	LEU
1	A	176	LEU
1	A	187	LEU
1	A	214	GLN
1	A	220	LEU
1	A	268	THR
1	A	277	ARG
1	A	293	ARG
1	A	301	LEU
1	A	314	ILE
1	A	317	ASN
1	A	318	HIS
1	A	320	LYS
1	A	323	ILE
1	A	324	ARG
1	A	337	LEU
1	A	359	LEU
1	A	388	LEU
1	A	390	TYR
1	A	443	CYS
1	A	456	THR
1	A	473	ILE
1	A	503	ARG
1	A	505	ILE
1	A	512	LEU
1	A	514	ASP
1	A	518	LEU
1	A	520	ARG

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Mol	Chain	Res	Type
1	A	530	LEU
1	A	538	THR
1	A	594	LYS
1	A	606	MET
1	A	628	MET
1	A	639	GLN
1	A	641	VAL
1	A	673	LEU
1	A	679	GLU
1	A	696	ILE
1	A	712	LEU
1	A	724	LEU
1	A	743	LEU
1	A	745	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	85	ASN
1	A	89	GLN
1	A	127	ASN
1	A	168	GLN
1	A	214	GLN
1	A	251	HIS
1	A	266	ASN
1	A	270	ASN
1	A	317	ASN
1	A	534	GLN
1	A	561	GLN
1	A	613	GLN
1	A	692	GLN
1	A	710	HIS
1	A	713	ASN

### 5.3.3 RNA ⓘ

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](#)

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	DGT	A	1001	2	26,33,33	0.98	1 (3%)	28,52,52	1.53	4 (14%)
4	ADP	A	1002	-	25,29,29	1.09	2 (8%)	24,45,45	1.75	2 (8%)

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	DGT	A	1001	2	-	0/18/34/34	0/3/3/3
4	ADP	A	1002	-	-	0/12/32/32	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1002	ADP	PB-O3A	2.09	1.63	1.60
4	A	1002	ADP	C5-C4	3.22	1.47	1.40
3	A	1001	DGT	C6-N1	3.24	1.38	1.33

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
4	A	1002	ADP	N3-C2-N1	-6.56	123.14	128.86
3	A	1001	DGT	N3-C2-N1	-4.92	120.27	127.46
4	A	1002	ADP	C4-C5-N7	-3.27	106.25	109.41
3	A	1001	DGT	C5-C6-N1	-2.25	120.28	123.48
3	A	1001	DGT	C6-N1-C2	2.50	119.65	116.06
3	A	1001	DGT	C2-N3-C4	4.30	120.17	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1001	DGT	1	0
4	A	1002	ADP	1	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	664/888 (74%)	0.34	50 (7%) 15 14	22, 38, 76, 98	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	145	TYR	7.6
1	A	665	THR	7.4
1	A	458	GLU	7.3
1	A	664	ILE	6.7
1	A	76	THR	6.5
1	A	638	PHE	5.9
1	A	459	ASP	5.8
1	A	77	LEU	5.5
1	A	81	ILE	5.4
1	A	146	PHE	5.2
1	A	461	LYS	5.1
1	A	78	ALA	4.9
1	A	390	TYR	4.4
1	A	628	MET	4.4
1	A	457	SER	4.3
1	A	89	GLN	4.3
1	A	630	SER	4.0
1	A	639	GLN	4.0
1	A	456	THR	3.9
1	A	79	ALA	3.9
1	A	462	THR	3.7
1	A	667	ASN	3.6
1	A	80	ARG	3.6
1	A	319	GLY	3.5
1	A	656	ASP	3.5
1	A	717	ARG	3.4
1	A	746	GLN	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	647	ARG	3.2
1	A	489	TYR	3.2
1	A	88	LYS	3.2
1	A	654	ILE	3.1
1	A	159	ILE	3.0
1	A	153	ARG	3.0
1	A	658	GLY	2.9
1	A	629	TYR	2.9
1	A	661	GLN	2.8
1	A	723	LYS	2.8
1	A	662	TYR	2.7
1	A	657	GLU	2.5
1	A	509	VAL	2.4
1	A	675	ASN	2.3
1	A	653	GLY	2.3
1	A	659	MET	2.3
1	A	92	LYS	2.2
1	A	446	ALA	2.2
1	A	85	ASN	2.2
1	A	149	LYS	2.2
1	A	317	ASN	2.1
1	A	82	ALA	2.0
1	A	90	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 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	ADP	A	1002	27/27	0.97	0.11	-0.89	36,40,41,43	0
3	DGT	A	1001	31/31	0.98	0.08	-1.12	21,27,30,33	0
2	MG	A	2001	1/1	0.92	0.29	-	43,43,43,43	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.