



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:33 PM UTC

PDB ID : 1XJJ / pdb_00001xjj
Title : Structural mechanism of allosteric substrate specificity in a ribonucleotide reductase: dGTP complex
Authors : Larsson, K.-M.; Jordan, A.; Eliasson, R.; Reichard, P.; Logan, D.T.; Nordlund, P.
Deposited on : 2004-09-23
Resolution : 1.86 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	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

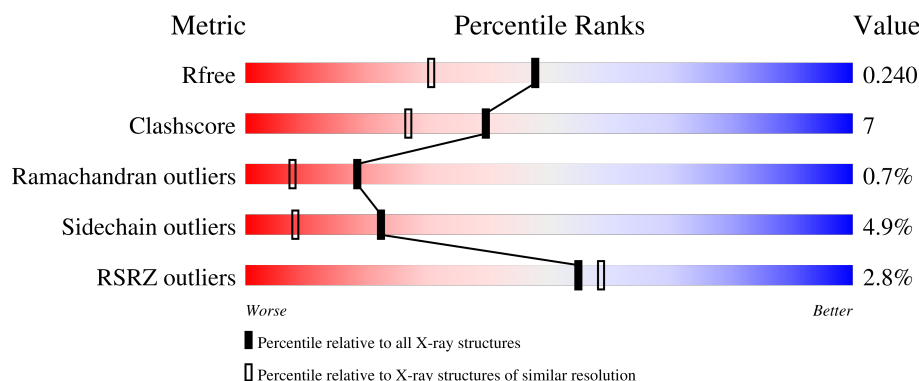
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

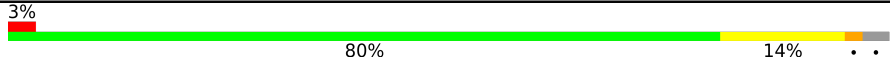
The reported resolution of this entry is 1.86 Å.

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	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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	644	
1	B	644	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10591 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 ribonucleotide reductase, B12-dependent.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	623	Total	C	N	O	S	0	0	0
			5005	3208	848	929	20			
1	B	611	Total	C	N	O	S	0	0	0
			4910	3149	833	908	20			

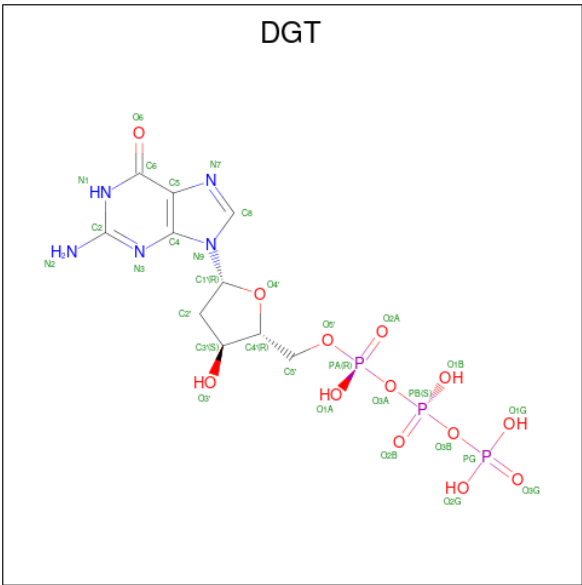
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	205	SER	TYR	SEE REMARK 999	UNP O33839
B	205	SER	TYR	SEE REMARK 999	UNP O33839

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 3 is 2'-DEOXYGUANOSINE-5'-TRIPHOSPHATE (CCD ID: DGT) (formula: C₁₀H₁₆N₅O₁₃P₃).



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

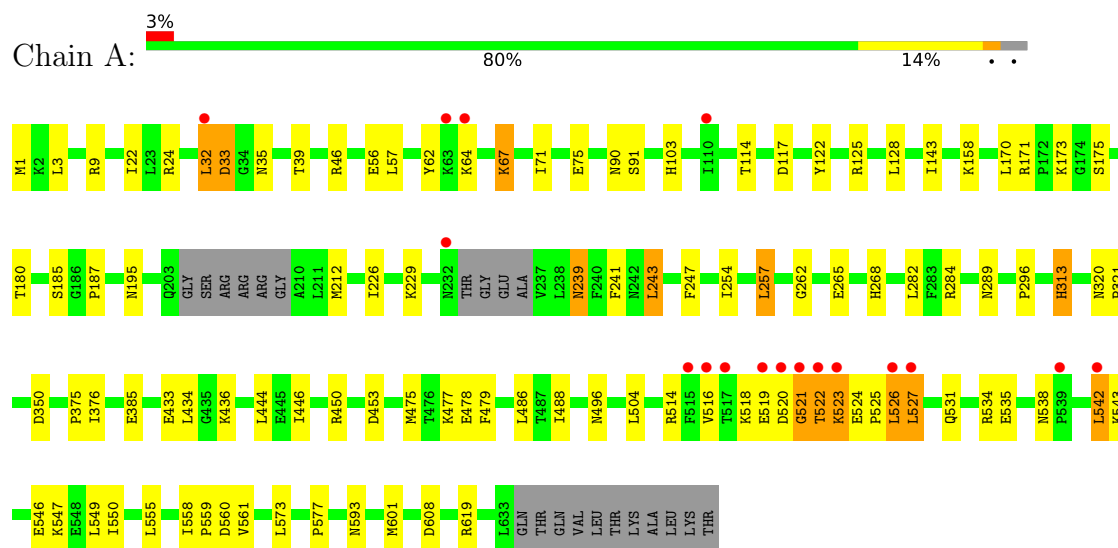
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	292	Total	O	0	0
			292	292		
4	B	258	Total	O	0	0
			258	258		

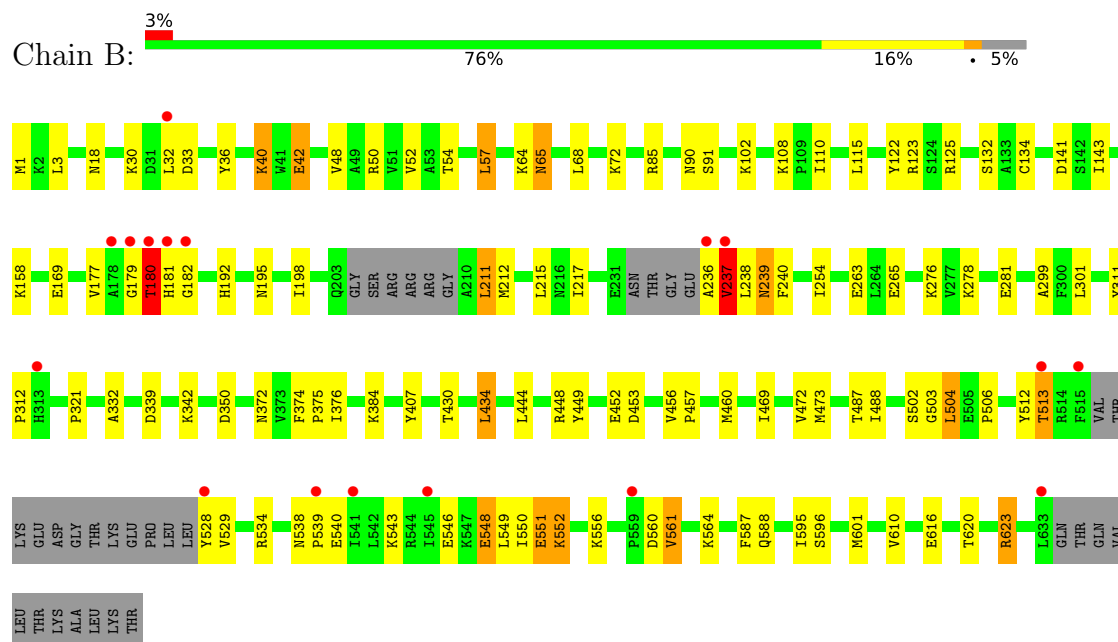
3 Residue-property plots [i](#)

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: ribonucleotide reductase, B12-dependent



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4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	118.07Å 123.88Å 106.30Å 90.00° 103.70° 90.00°	Depositor
Resolution (Å)	42.26 – 1.86 42.26 – 1.86	Depositor EDS
% Data completeness (in resolution range)	99.9 (42.26-1.86) 99.9 (42.26-1.86)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.21 (at 1.86Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.196 , 0.241 0.195 , 0.240	Depositor DCC
R_{free} test set	6217 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	25.4	Xtriage
Anisotropy	0.195	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 34.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10591	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.72% 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: DGT, MG

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	1.05	2/5104 (0.0%)	1.11	6/6891 (0.1%)
1	B	1.02	6/5007 (0.1%)	1.11	12/6758 (0.2%)
All	All	1.04	8/10111 (0.1%)	1.11	18/13649 (0.1%)

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	B	0	2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	561	VAL	CA-CB	7.55	1.60	1.54
1	B	587	PHE	C-O	-6.89	1.15	1.24
1	B	180	THR	CA-CB	6.50	1.64	1.53
1	B	299	ALA	CA-CB	5.84	1.61	1.53
1	A	247	PHE	C-O	-5.56	1.18	1.24
1	A	488	ILE	CA-CB	5.30	1.61	1.54
1	B	211	LEU	C-O	-5.07	1.17	1.23
1	B	217	ILE	CA-CB	5.03	1.61	1.54

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	374	PHE	CA-C-N	7.53	127.70	119.87
1	B	374	PHE	C-N-CA	7.53	127.70	119.87
1	A	91	SER	CA-C-N	-7.05	112.37	119.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	91	SER	C-N-CA	-7.05	112.37	119.56
1	B	301	LEU	N-CA-C	6.30	117.81	111.07
1	B	623	ARG	NE-CZ-NH1	5.98	127.48	121.50
1	A	573	LEU	N-CA-C	5.82	120.23	113.18
1	A	67	LYS	N-CA-C	5.79	118.06	111.11
1	B	623	ARG	CB-CG-CD	5.76	124.54	111.30
1	B	237	VAL	CA-C-N	5.69	131.94	121.70
1	B	237	VAL	C-N-CA	5.69	131.94	121.70
1	B	538	ASN	CA-C-N	-5.57	113.96	119.92
1	B	538	ASN	C-N-CA	-5.57	113.96	119.92
1	B	623	ARG	NE-CZ-NH2	-5.54	114.21	119.20
1	B	91	SER	N-CA-C	5.54	121.08	113.77
1	A	268	HIS	CA-C-N	-5.39	114.11	119.56
1	A	268	HIS	C-N-CA	-5.39	114.11	119.56
1	B	180	THR	N-CA-C	-5.22	99.68	110.80

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	237	VAL	Peptide
1	B	560	ASP	Peptide

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	5005	0	5057	63	0
1	B	4910	0	4958	78	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	62	0	24	0	0
3	B	62	0	24	0	0
4	A	292	0	0	5	0
4	B	258	0	0	3	0
All	All	10591	0	10063	135	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 7.

All (135) 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:522:THR:HB	1:A:523:LYS:HA	1.41	1.02
1:B:449:TYR:HE1	1:B:473:MET:HE1	1.22	0.99
1:B:601:MET:HE1	1:B:610:VAL:HG22	1.44	0.96
1:A:128:LEU:HD13	1:B:179:GLY:HA2	1.47	0.95
1:B:180:THR:HG22	1:B:181:HIS:H	1.30	0.94
1:B:487:THR:O	1:B:488:ILE:HD13	1.64	0.94
1:A:128:LEU:CD1	1:B:179:GLY:HA2	2.00	0.92
1:B:449:TYR:HE1	1:B:473:MET:CE	1.83	0.91
1:B:195:ASN:HD21	1:B:240:PHE:H	1.14	0.87
1:B:449:TYR:CE1	1:B:473:MET:HE1	2.10	0.86
1:B:457:PRO:HD3	1:B:473:MET:HE2	1.58	0.85
1:B:551:GLU:O	1:B:552:LYS:HB2	1.78	0.81
1:B:512:TYR:O	1:B:513:THR:HG23	1.81	0.80
1:A:542:LEU:O	1:A:546:GLU:HG3	1.87	0.75
1:A:385:GLU:OE1	4:A:1262:HOH:O	2.05	0.75
1:B:64:LYS:O	1:B:65:ASN:HB2	1.88	0.72
1:A:522:THR:CB	1:A:523:LYS:HA	2.18	0.72
1:B:1:MET:HB3	1:B:350:ASP:OD2	1.89	0.72
1:B:239:ASN:H	1:B:239:ASN:HD22	1.37	0.71
1:B:180:THR:HG22	1:B:181:HIS:N	2.04	0.70
1:A:524:GLU:HG2	1:A:525:PRO:HD2	1.74	0.70
1:A:619:ARG:NH1	4:A:1204:HOH:O	2.26	0.69
1:A:241:PHE:HB3	1:A:243:LEU:HD13	1.75	0.69
1:B:457:PRO:HD3	1:B:473:MET:CE	2.24	0.68
1:A:1:MET:HB3	1:A:350:ASP:OD2	1.96	0.66
1:B:180:THR:CG2	1:B:181:HIS:H	2.09	0.65
1:B:623:ARG:HD2	4:B:1017:HOH:O	1.97	0.64
1:B:263:GLU:HG2	1:B:278:LYS:HD3	1.80	0.64
1:A:475:MET:O	1:A:478:GLU:HG2	1.98	0.64
1:B:449:TYR:CE1	1:B:473:MET:CE	2.74	0.63
1:B:122:TYR:O	1:B:125:ARG:HD3	1.99	0.63
1:B:212:MET:HB2	1:B:321:PRO:HA	1.81	0.62
1:B:192:HIS:HE1	1:B:236:ALA:O	1.82	0.62
1:A:24:ARG:HH22	1:A:39:THR:HB	1.64	0.62
1:A:284:ARG:NH2	1:A:608:ASP:OD1	2.27	0.59
1:B:18:ASN:ND2	1:B:528:TYR:OH	2.35	0.59
1:A:549:LEU:HD13	1:A:555:LEU:HD23	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:226:ILE:HG22	1:A:289:ASN:HD22	1.67	0.58
1:B:539:PRO:O	1:B:540:GLU:HB2	2.03	0.58
1:A:257:LEU:HD22	1:A:262:GLY:HA3	1.85	0.58
1:B:64:LYS:HE3	1:B:115:LEU:HD21	1.86	0.58
1:A:577:PRO:HG3	1:A:601:MET:HG2	1.86	0.57
1:A:128:LEU:HD11	1:B:179:GLY:HA2	1.84	0.57
1:A:518:LYS:H	1:A:522:THR:CG2	2.18	0.57
1:B:40:LYS:HB3	1:B:40:LYS:NZ	2.20	0.57
1:B:265:GLU:OE1	1:B:276:LYS:HG2	2.05	0.56
1:A:195:ASN:HD21	1:A:239:ASN:ND2	2.04	0.55
1:B:180:THR:CG2	1:B:181:HIS:N	2.67	0.55
1:B:548:GLU:O	1:B:551:GLU:O	2.23	0.55
1:B:40:LYS:NZ	1:B:40:LYS:CB	2.70	0.55
1:A:521:GLY:HA3	1:A:522:THR:O	2.06	0.54
1:A:212:MET:HB2	1:A:321:PRO:HA	1.89	0.54
1:B:50:ARG:NH1	1:B:108:LYS:O	2.37	0.54
1:A:229:LYS:HG3	1:A:243:LEU:HD22	1.90	0.54
1:A:313:HIS:HD2	1:A:446:ILE:HD12	1.73	0.53
1:B:40:LYS:HB3	1:B:40:LYS:HZ3	1.73	0.53
1:B:64:LYS:O	1:B:65:ASN:CB	2.57	0.53
1:A:523:LYS:HB3	1:A:524:GLU:HA	1.90	0.52
1:A:239:ASN:HD22	1:A:239:ASN:H	1.57	0.52
1:B:195:ASN:ND2	1:B:240:PHE:H	1.96	0.52
1:B:456:VAL:HG11	1:B:460:MET:HE2	1.93	0.50
1:A:516:VAL:HG23	1:A:526:LEU:HD22	1.94	0.50
1:A:103:HIS:HE1	4:A:1191:HOH:O	1.94	0.50
1:B:134:CYS:O	1:B:332:ALA:HA	2.12	0.50
1:B:488:ILE:HG21	1:B:504:LEU:HG	1.93	0.50
1:A:453:ASP:HA	4:A:1233:HOH:O	2.13	0.49
1:A:375:PRO:C	1:A:376:ILE:HG13	2.38	0.49
1:A:433:GLU:CA	1:A:436:LYS:HE3	2.43	0.49
1:B:453:ASP:HA	4:B:1211:HOH:O	2.12	0.49
1:A:128:LEU:HD13	1:B:179:GLY:CA	2.32	0.49
1:B:469:ILE:O	1:B:473:MET:HG2	2.14	0.48
1:B:239:ASN:HD22	1:B:239:ASN:N	2.09	0.48
1:B:430:THR:HG22	1:B:434:LEU:HD22	1.95	0.48
1:A:143:ILE:HG13	1:B:158:LYS:HD2	1.95	0.48
1:B:50:ARG:NH2	1:B:108:LYS:O	2.46	0.48
1:A:433:GLU:HA	1:A:436:LYS:HE3	1.95	0.47
1:A:46:ARG:HH21	1:A:75:GLU:HG2	1.78	0.47
1:A:56:GLU:HB2	1:A:71:ILE:HG12	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:30:LYS:HE2	1:B:36:TYR:CE1	2.49	0.47
1:B:529:VAL:HB	1:B:534:ARG:NH2	2.30	0.47
1:B:448:ARG:O	1:B:452:GLU:HB2	2.15	0.46
1:A:313:HIS:CD2	1:A:446:ILE:HD12	2.51	0.46
1:A:122:TYR:O	1:A:125:ARG:HD3	2.15	0.46
1:B:311:TYR:CD1	1:B:312:PRO:HA	2.51	0.46
1:A:531:GLN:HE21	1:A:534:ARG:HH21	1.62	0.46
1:A:62:TYR:HB2	1:A:67:LYS:HE3	1.97	0.46
1:B:601:MET:HE1	1:B:610:VAL:CG2	2.31	0.46
1:B:588:GLN:CD	1:B:623:ARG:HD3	2.41	0.45
1:B:54:THR:O	1:B:57:LEU:HB2	2.15	0.45
1:B:339:ASP:OD1	1:B:339:ASP:C	2.59	0.45
1:A:22:ILE:CD1	1:A:496:ASN:HB3	2.47	0.45
1:B:68:LEU:O	1:B:72:LYS:HG3	2.17	0.45
1:A:320:ASN:HB2	1:A:321:PRO:HD2	1.99	0.44
1:B:502:SER:O	1:B:503:GLY:C	2.59	0.44
1:B:265:GLU:OE2	1:B:276:LYS:HE3	2.17	0.44
1:A:170:LEU:HD12	1:A:187:PRO:HA	1.99	0.44
1:A:450:ARG:HA	1:A:477:LYS:HG3	1.99	0.44
1:B:588:GLN:HB2	1:B:595:ILE:HD12	1.99	0.44
1:A:33:ASP:HB2	1:A:35:ASN:HD22	1.83	0.43
1:B:72:LYS:O	1:B:110:ILE:HD11	2.18	0.43
1:B:237:VAL:HA	1:B:238:LEU:HB2	1.99	0.43
1:A:478:GLU:HG3	1:A:479:PHE:CD2	2.54	0.43
1:A:527:LEU:CD1	1:A:550:ILE:HD11	2.47	0.43
1:A:114:THR:O	1:A:117:ASP:HB2	2.18	0.43
1:A:527:LEU:HD12	1:A:550:ILE:HD11	1.99	0.43
1:B:616:GLU:O	1:B:620:THR:HG23	2.19	0.43
1:A:226:ILE:HG22	1:A:289:ASN:ND2	2.32	0.43
1:B:375:PRO:C	1:B:376:ILE:HG13	2.43	0.43
1:B:85:ARG:O	1:B:342:LYS:NZ	2.52	0.42
1:B:181:HIS:HB3	1:B:182:GLY:HA2	2.00	0.42
1:A:9:ARG:NH2	4:A:1095:HOH:O	2.52	0.42
1:B:384:LYS:HG3	4:B:1123:HOH:O	2.19	0.42
1:A:486:LEU:O	1:A:593:ASN:HB2	2.20	0.42
1:A:523:LYS:H	1:A:524:GLU:CB	2.32	0.42
1:A:313:HIS:N	1:A:313:HIS:ND1	2.68	0.41
1:B:215:LEU:HA	1:B:215:LEU:HD12	1.90	0.41
1:B:546:GLU:O	1:B:550:ILE:HG13	2.19	0.41
1:A:32:LEU:H	1:A:32:LEU:HG	1.69	0.41
1:B:177:VAL:C	1:B:179:GLY:H	2.28	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:469:ILE:O	1:B:472:VAL:HG12	2.20	0.41
1:A:173:LYS:HG3	1:A:185:SER:O	2.20	0.41
1:A:518:LYS:H	1:A:522:THR:HG22	1.85	0.41
1:B:40:LYS:HG3	1:B:42:GLU:HG2	2.02	0.41
1:B:198:ILE:HG13	1:B:211:LEU:HD11	2.03	0.41
1:A:158:LYS:HD2	1:B:143:ILE:HG13	2.03	0.41
1:A:171:ARG:HG3	1:A:175:SER:HB2	2.02	0.41
1:A:518:LYS:O	1:A:522:THR:HG23	2.21	0.41
1:B:212:MET:HB2	1:B:212:MET:HE2	1.94	0.41
1:A:22:ILE:HD13	1:A:496:ASN:HB3	2.03	0.41
1:A:558:ILE:HA	1:A:559:PRO:HD2	1.96	0.40
1:B:48:VAL:O	1:B:52:VAL:HG23	2.21	0.40
1:B:141:ASP:HB2	1:B:169:GLU:O	2.20	0.40
1:B:132:SER:OG	1:B:372:ASN:ND2	2.53	0.40
1:B:407:TYR:CZ	1:B:506:PRO:HD3	2.56	0.40
1:A:433:GLU:O	1:A:436:LYS:HE3	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	617/644 (96%)	595 (96%)	20 (3%)	2 (0%)	36	25
1	B	603/644 (94%)	574 (95%)	23 (4%)	6 (1%)	12	4
All	All	1220/1288 (95%)	1169 (96%)	43 (4%)	8 (1%)	18	8

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	65	ASN
1	B	180	THR

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Mol	Chain	Res	Type
1	B	32	LEU
1	B	513	THR
1	B	552	LYS
1	A	522	THR
1	B	556	LYS
1	A	521	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	550/566 (97%)	519 (94%)	31 (6%)	19 6
1	B	538/566 (95%)	516 (96%)	22 (4%)	27 12
All	All	1088/1132 (96%)	1035 (95%)	53 (5%)	22 8

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	32	LEU
1	A	33	ASP
1	A	57	LEU
1	A	64	LYS
1	A	90	ASN
1	A	180	THR
1	A	239	ASN
1	A	243	LEU
1	A	254	ILE
1	A	257	LEU
1	A	265	GLU
1	A	282	LEU
1	A	296	PRO
1	A	313	HIS
1	A	434	LEU
1	A	444	LEU

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Mol	Chain	Res	Type
1	A	504	LEU
1	A	514	ARG
1	A	519	GLU
1	A	520	ASP
1	A	523	LYS
1	A	526	LEU
1	A	527	LEU
1	A	535	GLU
1	A	538	ASN
1	A	542	LEU
1	A	543	LYS
1	A	547	LYS
1	A	560	ASP
1	A	561	VAL
1	B	3	LEU
1	B	33	ASP
1	B	40	LYS
1	B	42	GLU
1	B	57	LEU
1	B	90	ASN
1	B	102	LYS
1	B	123	ARG
1	B	180	THR
1	B	239	ASN
1	B	254	ILE
1	B	281	GLU
1	B	434	LEU
1	B	444	LEU
1	B	504	LEU
1	B	543	LYS
1	B	548	GLU
1	B	549	LEU
1	B	551	GLU
1	B	561	VAL
1	B	564	LYS
1	B	596	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	ASN
1	A	127	HIS

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Mol	Chain	Res	Type
1	A	181	HIS
1	A	239	ASN
1	A	289	ASN
1	A	345	ASN
1	A	346	ASN
1	A	418	ASN
1	A	464	ASN
1	A	507	ASN
1	A	531	GLN
1	A	603	GLN
1	B	18	ASN
1	B	96	ASN
1	B	127	HIS
1	B	192	HIS
1	B	195	ASN
1	B	239	ASN
1	B	345	ASN
1	B	346	ASN
1	B	359	GLN
1	B	464	ASN
1	B	507	ASN
1	B	603	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	B	1004	-	32,33,33	1.76	4 (12%)	48,52,52	2.02	15 (31%)
3	DGT	A	1001	2	32,33,33	2.27	7 (21%)	48,52,52	1.65	10 (20%)
3	DGT	B	1002	2	32,33,33	3.48	4 (12%)	48,52,52	1.74	10 (20%)
3	DGT	A	1003	-	32,33,33	1.92	4 (12%)	48,52,52	2.14	23 (47%)

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	B	1004	-	-	5/22/34/34	0/3/3/3
3	DGT	A	1001	2	-	2/22/34/34	0/3/3/3
3	DGT	B	1002	2	-	2/22/34/34	0/3/3/3
3	DGT	A	1003	-	-	5/22/34/34	0/3/3/3

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1002	DGT	PB-O3A	18.04	1.79	1.59
3	A	1001	DGT	PB-O3A	8.47	1.68	1.59
3	A	1003	DGT	PA-O3A	6.40	1.66	1.59
3	A	1003	DGT	PB-O3B	5.71	1.65	1.59
3	B	1004	DGT	PB-O3A	5.19	1.65	1.59
3	A	1001	DGT	PA-O3A	5.05	1.64	1.59
3	B	1004	DGT	PB-O3B	4.85	1.64	1.59
3	B	1002	DGT	PG-O3G	4.15	1.63	1.50
3	B	1004	DGT	PG-O3G	3.57	1.61	1.50
3	A	1001	DGT	PG-O3G	2.92	1.59	1.50
3	B	1002	DGT	C4-N9	-2.61	1.31	1.38
3	A	1001	DGT	C5-N7	-2.59	1.33	1.39
3	A	1001	DGT	PG-O1G	-2.53	1.45	1.54
3	B	1002	DGT	PB-O3B	2.41	1.62	1.59
3	A	1003	DGT	PG-O3G	2.40	1.57	1.50
3	B	1004	DGT	C5-N7	-2.38	1.34	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1001	DGT	C2-N1	2.29	1.43	1.37
3	A	1003	DGT	C5-N7	-2.12	1.34	1.39
3	A	1001	DGT	PB-O3B	-2.05	1.57	1.59

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1003	DGT	C2-N3-C4	5.36	121.53	112.30
3	B	1004	DGT	C2-N3-C4	5.34	121.49	112.30
3	B	1002	DGT	C2'-C1'-N9	-4.72	102.02	113.81
3	B	1002	DGT	C2-N3-C4	4.59	120.21	112.30
3	B	1004	DGT	O3A-PA-O2A	-4.55	97.01	110.70
3	B	1004	DGT	C5-C4-N3	-4.53	121.18	128.39
3	A	1001	DGT	C2-N3-C4	4.04	119.26	112.30
3	A	1003	DGT	C5-C4-N3	-4.04	121.95	128.39
3	A	1003	DGT	O4'-C4'-C5'	-4.03	96.43	109.33
3	A	1001	DGT	C5-C4-N3	-3.95	122.09	128.39
3	B	1004	DGT	O1A-PA-O3A	3.51	116.75	107.27
3	B	1004	DGT	O4'-C1'-N9	3.42	113.94	107.86
3	B	1004	DGT	N9-C4-N3	3.42	132.79	125.95
3	B	1002	DGT	O2G-PG-O1G	3.34	120.33	107.80
3	A	1003	DGT	O3A-PA-O2A	-3.27	100.87	110.70
3	A	1001	DGT	C2-N1-C6	-3.26	119.20	125.11
3	A	1001	DGT	O2G-PG-O1G	3.23	119.90	107.80
3	A	1003	DGT	C2-N1-C6	-3.23	119.26	125.11
3	B	1002	DGT	C5-C4-N3	-3.18	123.33	128.39
3	B	1004	DGT	C2-N1-C6	-3.16	119.38	125.11
3	B	1004	DGT	C4'-O4'-C1'	-3.12	102.11	109.51
3	A	1003	DGT	O2G-PG-O3B	3.09	115.01	104.64
3	A	1003	DGT	N9-C8-N7	-3.02	107.81	113.40
3	A	1001	DGT	N9-C4-N3	2.86	131.67	125.95
3	B	1004	DGT	O2G-PG-O1G	2.85	118.47	107.80
3	A	1003	DGT	O4'-C1'-N9	-2.81	102.88	107.86
3	A	1003	DGT	N9-C4-N3	2.81	131.56	125.95
3	A	1001	DGT	O1B-PB-O3A	-2.77	99.79	107.27
3	B	1002	DGT	O4'-C1'-C2'	-2.75	101.10	106.25
3	A	1003	DGT	O2G-PG-O1G	2.67	117.80	107.80
3	A	1001	DGT	O6-C6-C5	-2.62	119.62	126.53
3	A	1003	DGT	O1A-PA-O3A	2.60	114.29	107.27
3	A	1003	DGT	C8-N7-C5	2.56	108.81	104.26
3	B	1002	DGT	N9-C8-N7	-2.48	108.80	113.40
3	A	1003	DGT	C5-C6-N1	2.48	119.57	113.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1002	DGT	O6-C6-N1	-2.46	115.49	120.11
3	A	1003	DGT	O3B-PG-O3G	-2.44	98.20	111.04
3	B	1002	DGT	C2-N1-C6	-2.43	120.70	125.11
3	A	1001	DGT	C2'-C1'-N9	-2.42	107.75	113.81
3	A	1003	DGT	O3B-PB-O2B	-2.39	103.53	110.70
3	A	1003	DGT	C6-C5-N7	2.36	134.59	130.29
3	B	1002	DGT	C6-C5-N7	2.35	134.57	130.29
3	A	1003	DGT	C4'-O4'-C1'	-2.29	104.08	109.51
3	A	1003	DGT	O6-C6-C5	-2.28	120.51	126.53
3	B	1004	DGT	N9-C8-N7	-2.28	109.18	113.40
3	A	1001	DGT	C5-C6-N1	2.22	118.91	113.25
3	B	1004	DGT	O4'-C1'-C2'	-2.20	102.14	106.25
3	B	1004	DGT	C5-C6-N1	2.19	118.82	113.25
3	A	1003	DGT	C4-C5-N7	-2.18	107.21	110.67
3	B	1004	DGT	O6-C6-C5	-2.15	120.86	126.53
3	B	1004	DGT	O1A-PA-O2A	2.15	122.43	112.44
3	A	1003	DGT	O1B-PB-O3A	-2.14	101.50	107.27
3	A	1003	DGT	O1B-PB-O3B	2.12	113.01	107.27
3	A	1003	DGT	N1-C2-N3	-2.10	119.47	123.32
3	A	1003	DGT	O1B-PB-O2B	2.10	122.20	112.44
3	A	1001	DGT	O5'-C5'-C4'	2.06	116.00	108.99
3	B	1004	DGT	C8-N7-C5	2.05	107.92	104.26
3	B	1002	DGT	C5-C6-N1	2.01	118.38	113.25

There are no chirality outliers.

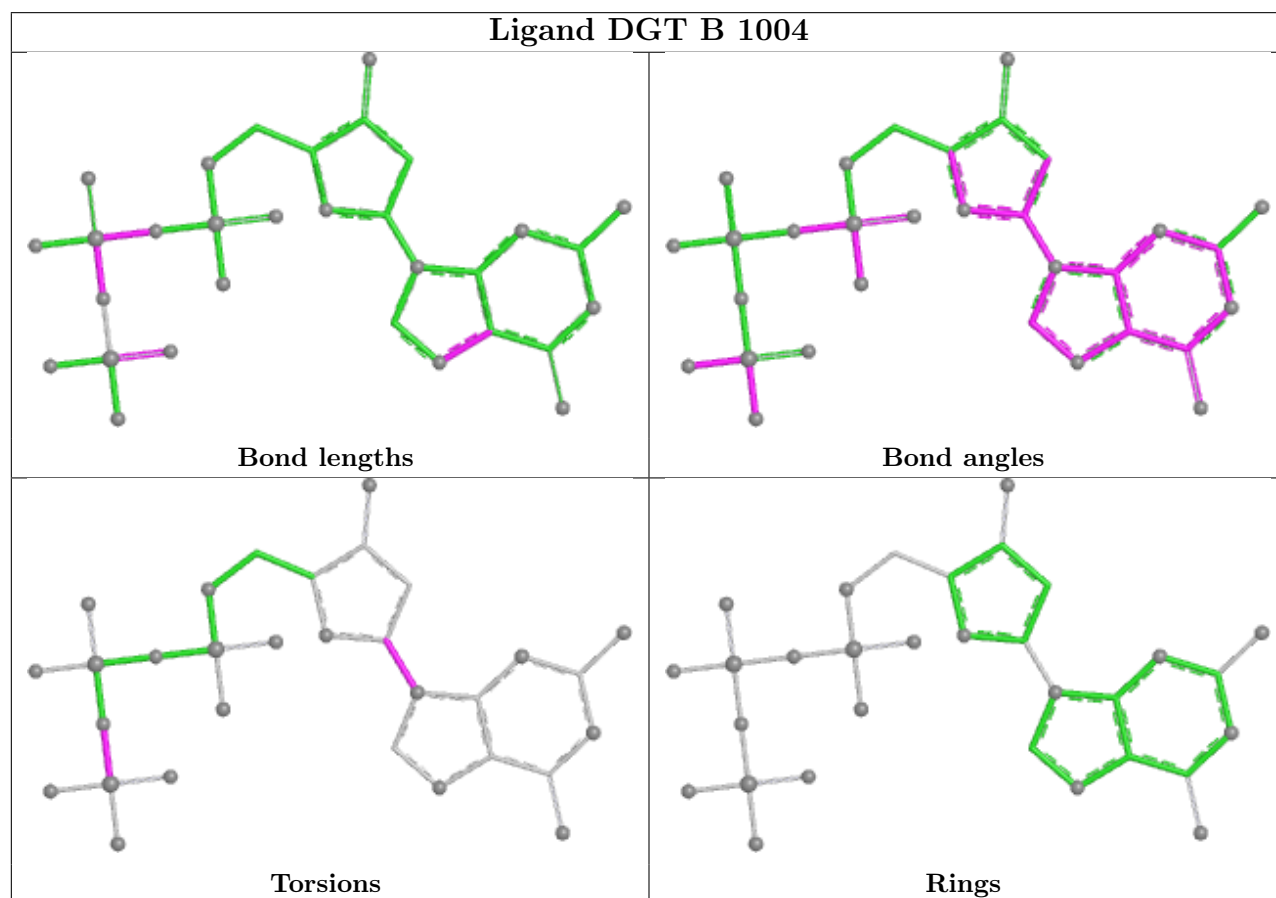
All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1003	DGT	PB-O3B-PG-O2G
3	B	1004	DGT	PB-O3B-PG-O1G
3	B	1004	DGT	PB-O3B-PG-O2G
3	B	1004	DGT	C2'-C1'-N9-C8
3	B	1004	DGT	C2'-C1'-N9-C4
3	A	1001	DGT	PB-O3B-PG-O2G
3	B	1002	DGT	PB-O3B-PG-O2G
3	A	1001	DGT	PA-O3A-PB-O2B
3	A	1003	DGT	PB-O3A-PA-O2A
3	B	1004	DGT	PB-O3B-PG-O3G
3	A	1003	DGT	PB-O3B-PG-O1G
3	A	1003	DGT	PB-O3B-PG-O3G
3	A	1003	DGT	PB-O3A-PA-O1A
3	B	1002	DGT	PA-O3A-PB-O1B

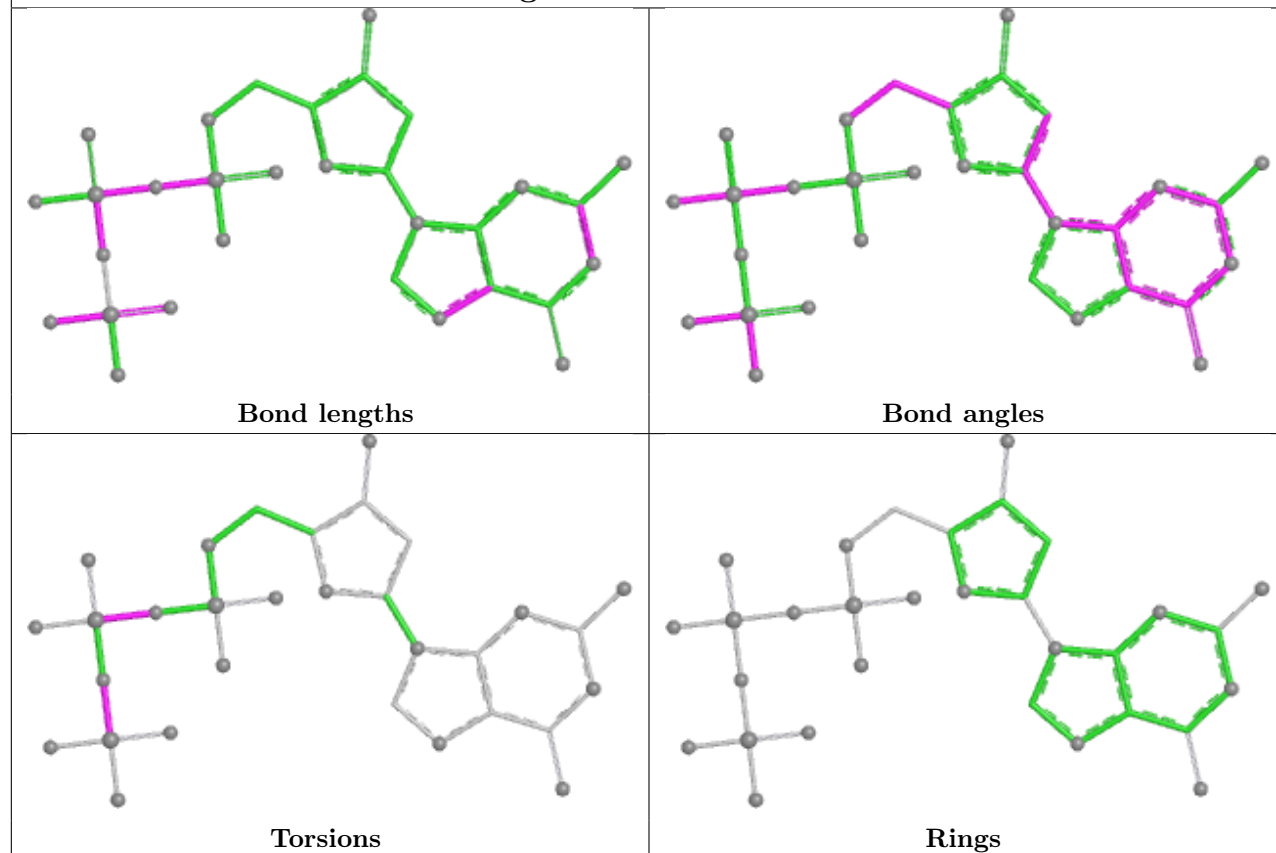
There are no ring outliers.

No monomer is involved in short contacts.

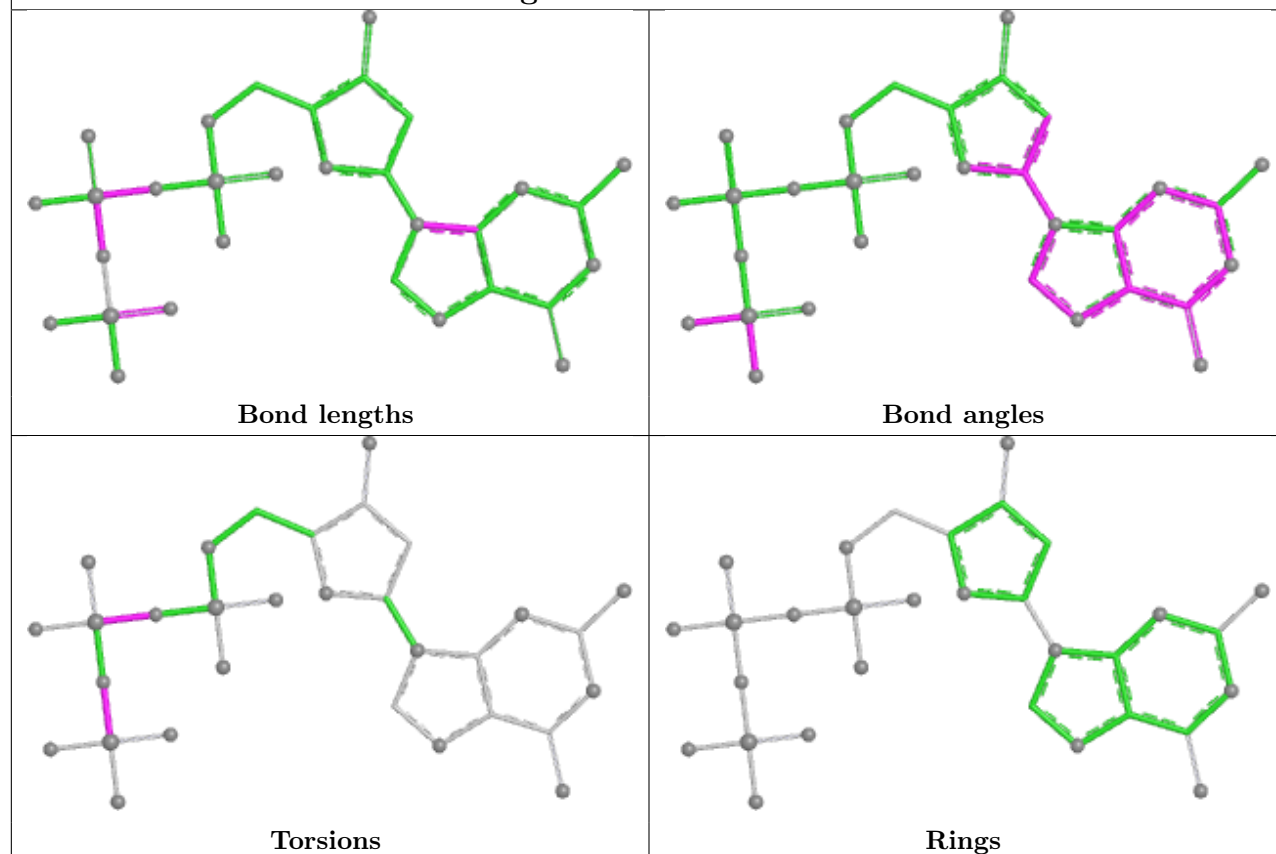
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

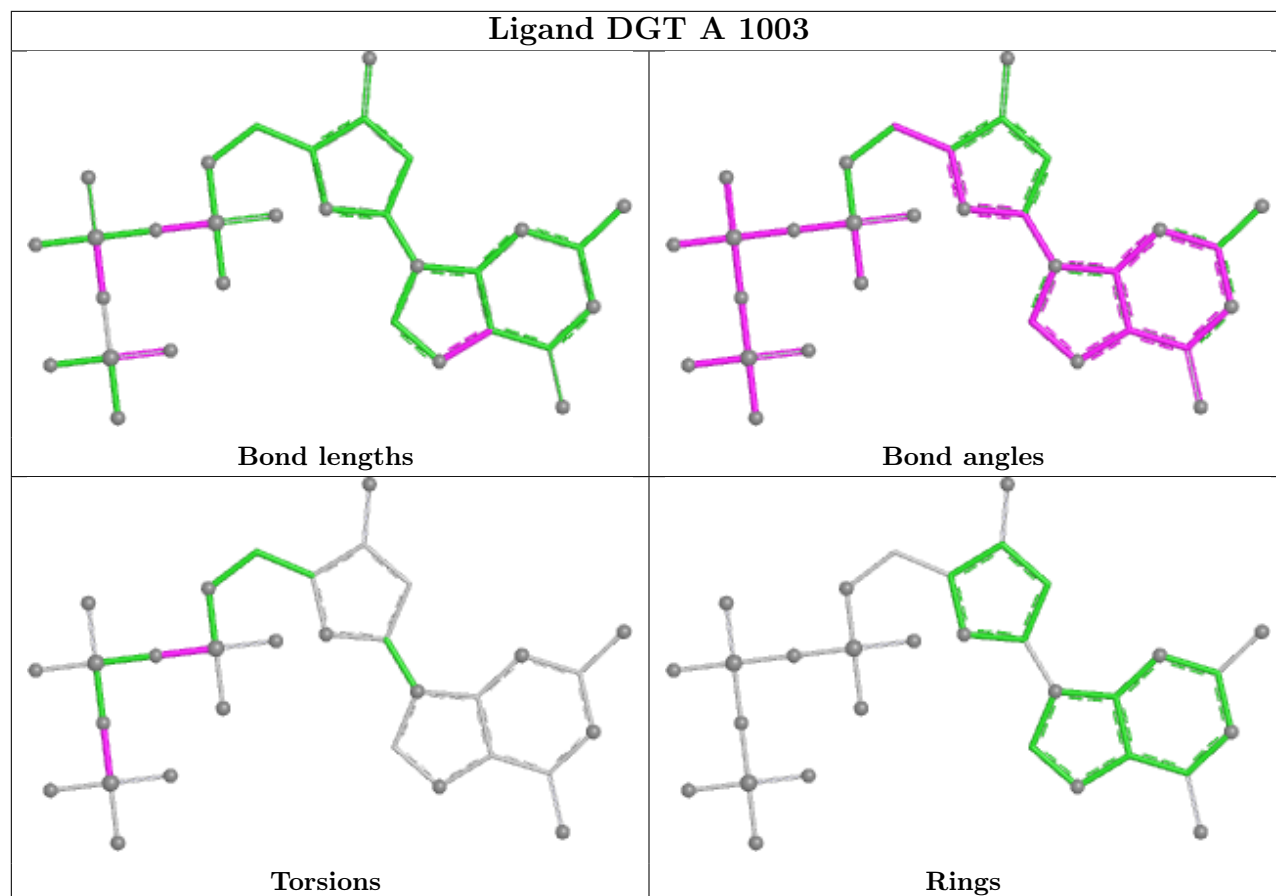


Ligand DGT A 1001



Ligand DGT B 1002





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	623/644 (96%)	0.05	17 (2%) 56 59	15, 27, 56, 88	0
1	B	611/644 (94%)	0.12	17 (2%) 55 58	17, 29, 59, 75	0
All	All	1234/1288 (95%)	0.08	34 (2%) 55 58	15, 28, 58, 88	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	236	ALA	4.4
1	B	180	THR	4.1
1	B	515	PHE	3.9
1	B	528	TYR	3.8
1	A	516	VAL	3.8
1	B	181	HIS	3.3
1	A	517	THR	3.0
1	A	521	GLY	3.0
1	A	526	LEU	2.9
1	B	633	LEU	2.9
1	A	522	THR	2.9
1	A	64	LYS	2.9
1	B	559	PRO	2.7
1	B	178	ALA	2.7
1	B	513	THR	2.7
1	B	545	ILE	2.7
1	A	515	PHE	2.6
1	B	539	PRO	2.6
1	B	182	GLY	2.6
1	B	313	HIS	2.6
1	B	179	GLY	2.6
1	B	237	VAL	2.5
1	A	523	LYS	2.4
1	B	541	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	232	ASN	2.3
1	B	32	LEU	2.3
1	A	520	ASP	2.2
1	A	519	GLU	2.2
1	A	539	PRO	2.2
1	A	110	ILE	2.2
1	A	32	LEU	2.1
1	A	527	LEU	2.1
1	A	63	LYS	2.1
1	A	542	LEU	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

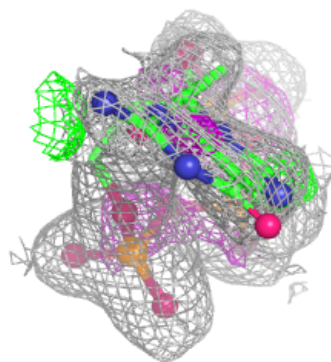
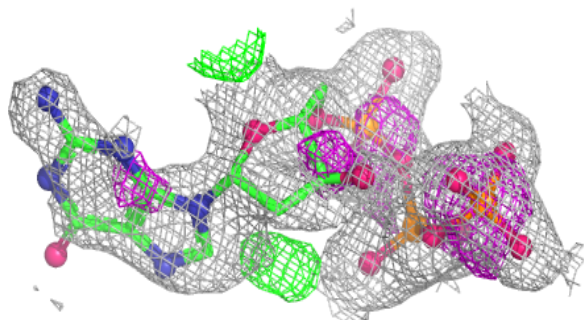
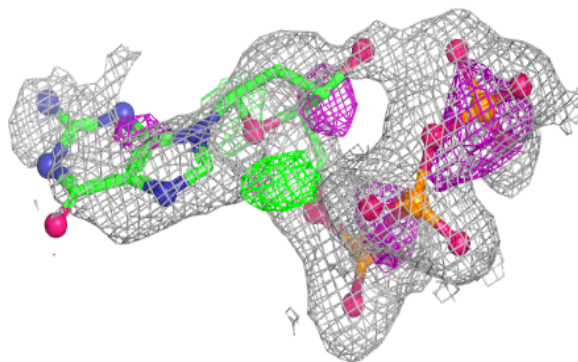
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	DGT	A	1003	31/31	0.93	0.13	26,63,75,75	0
3	DGT	B	1004	31/31	0.93	0.12	33,62,73,74	0
3	DGT	B	1002	31/31	0.96	0.06	14,21,44,44	0
3	DGT	A	1001	31/31	0.98	0.04	17,25,32,34	0
2	MG	B	1005	1/1	0.98	0.07	36,36,36,36	0
2	MG	A	1006	1/1	0.99	0.02	26,26,26,26	0

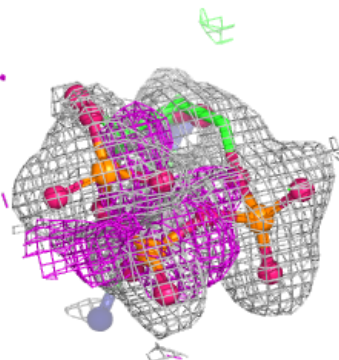
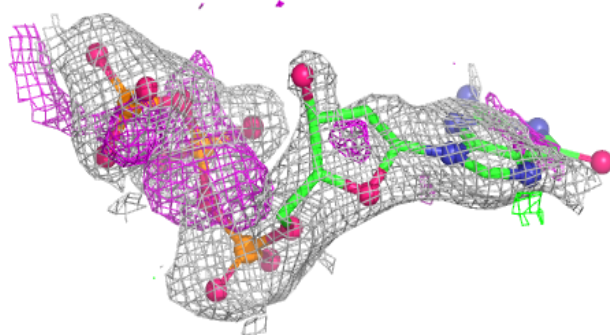
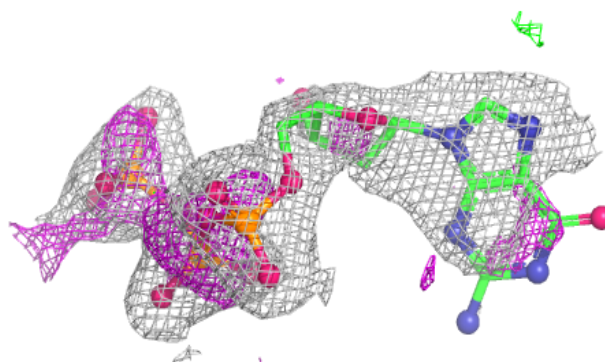
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around DGT A 1003:

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

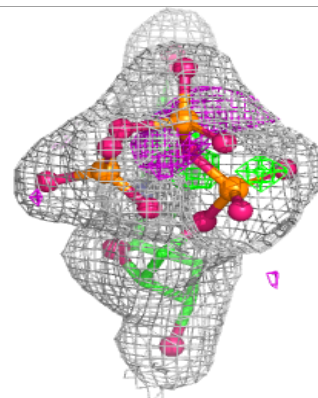
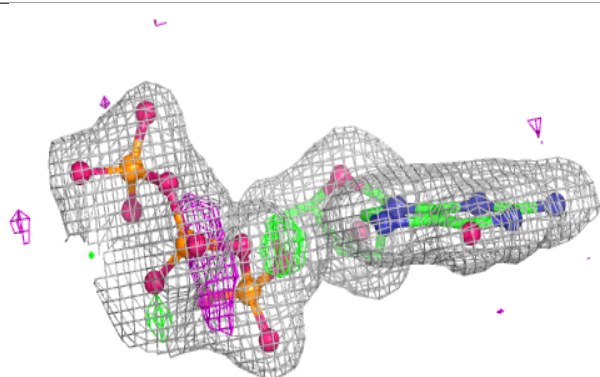
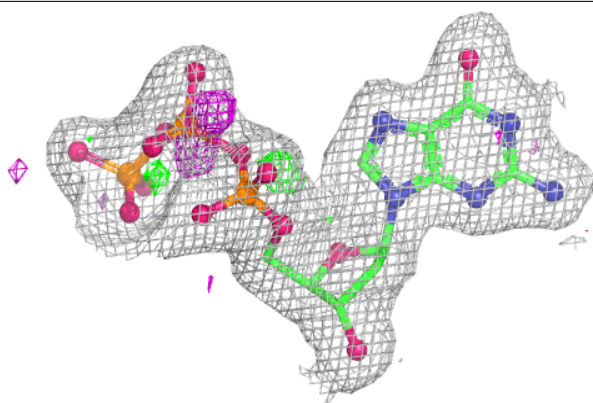
**Electron density around DGT B 1004:**

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

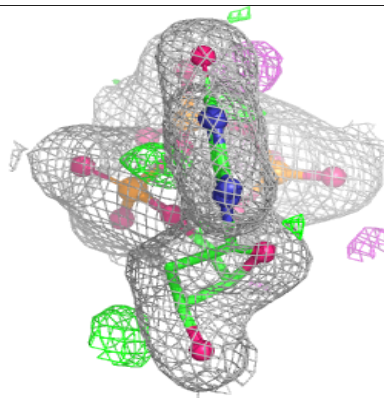
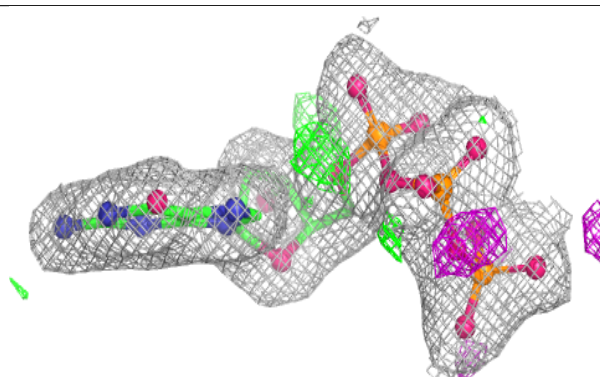
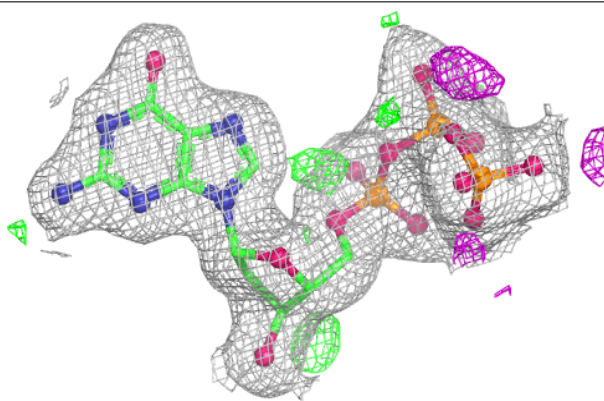


Electron density around DGT B 1002:

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

**Electron density around DGT 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 ⓘ

There are no such residues in this entry.